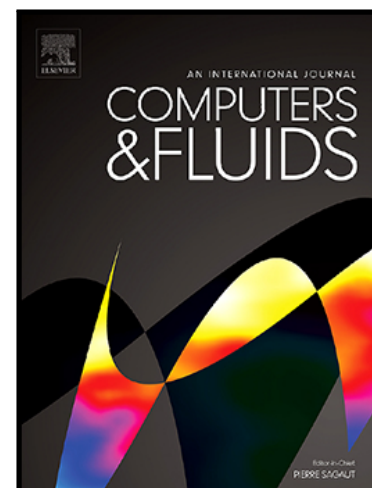


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## Highlights (for review)

### Highlights:

- CWENO/CWENOZ methods for mixed-element unstructured meshes
- Analysis of central stencil linear weight selection is given
- Tested for Euler and Navier-Stokes equations
- CWENO/CWENOZ methods more robust and efficient than classical WENO

# Arbitrary High Order Central Non-Oscillatory Schemes on Mixed-Element Unstructured Meshes

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## Abstract

In this paper we develop a family of very high-order central (up to 6th-order) non-oscillatory schemes for mixed-element unstructured meshes. The schemes are inherently compact in the sense that the central stencils employed are as compact as possible, and that the directional stencils are reduced in size therefore simplifying their implementation. Their key ingredient is the non-linear combination in a CWENO style similar to Dumbser et al [1] of a high-order polynomial arising from a central stencil with lower-order polynomials from directional stencils. Therefore, in smooth regions of the computational domain the optimum order of accuracy is recovered, while in regions of sharp-gradients the larger influence of the reconstructions from the directional stencils suppress the oscillations. It is the compactness of the directional stencils that increases the chances of at least one of them lying in a region with smooth data, that greatly enhances their robustness compared to classical WENO schemes. The two variants developed are CWENO and CWENOZ schemes, and it is the first time that such very-high-order schemes are designed for mixed-element unstructured meshes. We explore the influence of the linear weights in each of the schemes, and assess their performance in terms of accuracy, robustness and computational cost through a series of stringent 2D and 3D test problems. The results obtained demonstrate the improved robustness that the schemes offer, a parameter of paramount importance for and their potential use for industrial-scale engineering applications.

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## 1. Introduction

The development and application of high-order non-oscillatory schemes for computational fluid dynamics on unstructured meshes has been the focus of an overwhelming number of studies that spans many numerical frameworks and techniques. The catalyst for this development is firstly that arbitrary shaped unstructured elements can offer significant advantages for computational workflows that include complicated geometries that require rapid generated body fitted meshes for flows that are characterised by the presence of phenomena with contradicting requirements. On one hand flows regions with strong gradients or discontinuities require robust non-oscillatory schemes that can resolve the phenomena encountered in these regions without generating spurious oscillations, while on the other hand in regions of smooth flow features, such as vortices or acoustic waves, highly accurate schemes for resolving these delicate flow features is a necessity. Finding the perfect balance between these two contradicting requirements is the driving force upon which the saga towards the ultimate high-order non-oscillatory schemes has been based. To make this situation even more challenging this perfect balance must be obtained in a computationally efficient manner, something that the rapid changing high-performance computing (HPC) landscape makes it harder to sustain since the numerical schemes must be adapted or completely redesigned to make the most of the latest HPC hardware.

The plethora of numerical schemes available in this context are primarily developed in the Finite-Volume (FV) [1–16], discontinuous Galerkin (DG) [17–23], flux reconstruction (FR) [24–31] and spectral-finite-volume (SFV) [32, 33] frameworks. Although the frameworks are significantly different in principle the high-order spatial accuracy

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is achieved through a natural extension of the representation of the solution within each cell by polynomials. Two of the most popular families of schemes for providing non-oscillatory capabilities to a numerical framework is firstly the family of schemes that can detect when some solution bounds have been violated and switch to a lower-order approximation of the solution such as total-variation bounded (TVB) [34], total-variation diminishing (TVD) [35–37], monotone upstream scheme for conservation laws (MUSCL) [38–46], and multi-dimensional optimal order detection (MOOD) [3, 4, 47–49], and secondly by techniques where the spatial-regions with the best quality of information (smooth data) will have the largest influence in the approximation of the solution, such as the weighted essentially non-oscillatory (WENO) [2, 5, 6, 13, 50–55] schemes. It needs to be noted that high order WENO reconstruction in space is also the key ingredient of the ADER class of finite volume schemes of Toro and Titarev, see e.g. [56–59] and references therein for an overview. The first family is characterised by a compact support stencil (usually a fraction of the central stencil) which is beneficial for the computational efficiency, however these frameworks are not parameter-free and they are rather sensitive to the values of the parameters to define the behaviour of the algorithm once the bounds have been violated as previously reported in [38, 47]. On the other hand, the WENO family of methods offers a high-order accurate platform that is less sensitive to the parameters in particular if transformations are performed to reduce scaling effects [5, 6, 13], and a good computational efficiency, but the size of the directional stencils whose width is dependent on the quality of the mesh is increasing the complexity of the implementation of the schemes as well as their communication overhead across multiple-processes in an HPC setting. One of the ways to alleviate this is to reduce the size of the directional stencils and consequently the order of approximation from them. This strategy has been adopted by many in developing the sometimes called new, central, compact or cool WENO schemes [1, 12, 51, 53, 60–70]. The key benefit of using these schemes is the reduced computational cost compared to the traditional WENO schemes due to the reduced size of the directional stencils. The increased compactness of this class of schemes makes them more suitable to be deployed for frameworks such as DG, something that has already started as seen in [53, 69–71]. However, one of the most overlooked benefits of these type of CWENO schemes is an increased robustness, since the reduced size of the directional stencils translates directly into a higher probability of at least one of them lying in a region with smooth variation of data, compared to the traditional WENO schemes for unstructured meshes. The results obtained with these schemes show a larger sensitivity to the parameters involved, since the blending of a higher-order approximation emanating from the central stencil with lower-order approximation from the directional stencils requires a fine balance and calibration.

In the present study we are pursuing the extension of two variants of central WENO schemes to mixed-element unstructured meshes for an arbitrary high-order. The two variants developed are the CWENO and CWENOZ schemes, where in order to further reduce their computational footprint both variants use the stencil based compact (SBC) central stencil algorithm and Type 3 directional stencil algorithm developed in [72]. For the central stencil the polynomial of any desired order can be built, while for the directional stencils a  $\mathcal{P} = 1$  polynomial is employed requiring 4 and 6 neighbours in 2D and 3D respectively. The directional stencils are relatively compact and building them is straight forward since they are contained in the central stencil (at least for higher-order methods) and therefore making the implementation less complicated compared to the original WENO schemes. In this work we develop and apply schemes up to 6th-order of spatial accuracy for mixed-element meshes, and in principle they can be extended to any spatial order. Ultimately the goal is to make this class of very higher-order schemes more affordable and more robust for unstructured meshes even for industrial scale applications.

All the schemes are developed in the open source UCNS3D solver [73], and we assess their performance in terms of robustness, accuracy and computational efficiency for a series of stringent 2D and 3D test problems while also comparing them with the traditional WENO schemes. The paper is organized as follows. In Section 2 we introduce the numerical framework used to describe the high-order finite-volume framework for unstructured meshes, the reconstruction process for the WENO, CWENO and CWENOZ schemes, while describing the chosen fluxes and temporal discretisation employed. The numerical results obtained for all the test problems are presented in Section 3 and compared against analytical, reference or experimental solution whenever possible. Finally, the last section describes the conclusions drawn from this study.

## 2. Numerical Framework

Consider the unsteady non-linear hyperbolic system of conservation laws on a 3D domain  $\Omega$ , written in its conservative form:

$$\frac{\partial \mathbf{U}}{\partial t} + \nabla \cdot (\mathbf{F}(\mathbf{U})) = 0, \quad (1)$$

where  $\mathbf{U} = \mathbf{U}(\mathbf{x}, t)$  is the vector of conserved variables,  $\mathbf{x} = (x, y, z)$  denotes the coordinates of a point of the domain  $\Omega$ , and  $\mathbf{F}(\mathbf{U}) = (\mathbf{f}(\mathbf{U}), \mathbf{g}(\mathbf{U}), \mathbf{h}(\mathbf{U}))$  is the non-linear flux tensor. The physical domain  $\Omega$  consists of any combination of conforming tetrahedral, hexahedral, prism or pyramids in 3D, and quadrilateral or triangular 2D. All the elements are indexed by a unique mono-index  $i$ . Integrating Eq. (1) over the mesh element  $i$  using a high-order explicit finite-volume formulation the following equation is obtained:

$$\mathbf{U}_i^{n+1} = \mathbf{U}_i^n - \Delta t \frac{1}{|V_i|} \sum_{j=1}^{N_f} \sum_{\alpha=1}^{N_{qp}} F^{n_{ij}} \left( \mathbf{U}_{ij,L}^n(\mathbf{x}_{ij,\alpha}, t), \mathbf{U}_{ij,R}^n(\mathbf{x}_{ij,\alpha}, t) \right) \omega_\alpha |S_{ij}|, \quad (2)$$

where  $\mathbf{U}_i$  are the volume averaged conserved variables

$$\mathbf{U}_i = \frac{1}{|V_i|} \int_{V_i} \mathbf{U}(x, y, z) dV, \quad (3)$$

and  $F^{n_{ij}}$  is a numerical flux function in the direction normal to the cell interface between cell  $i$  and the neighbouring cell  $j$  as seen in Fig. 1,  $N_f$  is the number of faces per element,  $N_{qp}$  is the number of quadrature points used for approximating the surface integrals,  $|S_{ij}|$  is the surface area of the corresponding face, and  $\mathbf{U}_{ij,L}^n(\mathbf{x}_{ij,\alpha}, t)$  and  $\mathbf{U}_{ij,R}^n(\mathbf{x}_{ij,\alpha}, t)$  are the high-order approximations of the solutions for cell  $i$  and cell  $j$  respectively; while  $\alpha$  corresponds to different Gaussian integration points  $\mathbf{x}_\alpha$  and weights  $\omega_\alpha$  over each face. The volume, surface and line integrals are numerically approximated by suitable quadrature rules, see [74] for details on numerical approximations of multiple integrals.

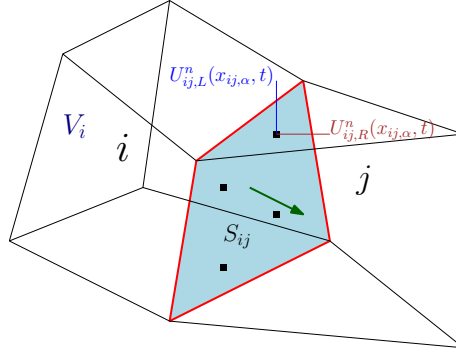


Figure 1: Drawing illustrating the interface between the considered cell  $i$  and its neighbour  $j$  highlighting the normal vector and the quadrature points at the interface.

### 2.1. Reconstruction

For a cell  $i$  a high-order polynomial  $p_i(x, y, z)$  of order  $r$  can be built that provides  $r + 1$  order of accuracy, by requiring it to have the same average as a general quantity  $\mathbf{U}_i$ . This can be formulated as:

$$\mathbf{U}_i = \frac{1}{|V_i|} \int_{V_i} p_i(x, y, z) dV. \quad (4)$$

The present polynomial reconstruction is based upon the approaches of [5, 6, 13, 52], that have been applied to smooth and discontinuous flow problems [38, 47, 72, 75–86] and only the key-components will be presented herein and the

reader is referred to [5, 6, 52] for further details. Unstructured meshes involve the presence of various elements shapes and sizes, and a transformation from physical space to a reference space is performed to reduce the scaling effects as introduced by Dumbser et al. [13, 14, 87]. In this study this is performed by decomposing each element into triangular or tetrahedral elements and using one of the decomposed elements as the reference element for transforming to the new system of coordinates as suggested by Tsoutsanis et al. [5, 6, 52]. The decomposition strategy is discussed in [6].

Let  $\mathbf{v}\mathbf{x}_{ij}$ ,  $j = 1, 2, \dots, J_i$  be the vertices of the considered 3D general element. Non tetrahedral elements are decomposed into tetrahedrals and one of them is chosen with  $\mathbf{w}_1 = (x_1, y_1, z_1)$ ,  $\mathbf{w}_2 = (x_2, y_2, z_2)$ ,  $\mathbf{w}_3 = (x_3, y_3, z_3)$  and  $\mathbf{w}_4 = (x_4, y_4, z_4)$  being its four vertices. The transformation from the Cartesian coordinates  $x, y, z$  into a reference space  $\xi, \eta, \zeta$  is given by the following equations:

$$\begin{pmatrix} x \\ y \\ z \end{pmatrix} = \begin{pmatrix} x_1 \\ y_1 \\ z_1 \end{pmatrix} + J \cdot \begin{pmatrix} \xi \\ \eta \\ \zeta \end{pmatrix}, \quad (5)$$

with the Jacobian matrix given by:

$$J = \begin{bmatrix} x_2 - x_1 & x_3 - x_1 & x_4 - x_1 \\ y_2 - y_1 & y_3 - y_1 & y_4 - y_1 \\ z_2 - z_1 & z_3 - z_1 & z_4 - z_1 \end{bmatrix}. \quad (6)$$

Using an inverse mapping the element  $V_i$  can be transformed to the element  $V'_i$  in the reference co-ordinate system as:

$$\mathbf{v}\mathbf{x}'_{ij} = J^{-1} \cdot (\mathbf{v}\mathbf{x}_{ij} - \mathbf{w}_1), \quad j = 1, 2, \dots, J_i. \quad (7)$$

and the spatial average of the conserved variable  $\mathbf{U}_i$  does not change during transformation

$$\mathbf{U}_i = \frac{1}{|V_i|} \int_{V_i} \mathbf{U}(x, y, z) dV \equiv \frac{1}{|V'_i|} \int_{V'_i} \mathbf{U}(\xi, \eta, \zeta) d\xi d\eta d\zeta. \quad (8)$$

The reconstruction is performed by building a central stencil  $\mathcal{S}^1$  by recursively adding neighbouring elements, consisting of  $M + 1$  cells including the considered cell  $i$ . In this study the stencil based compact algorithm (SBC) introduced in [72] is used due to its low computational cost and enhanced robustness, the reader is referred to [72] for more details. In the present study we employ  $M = 2K$ , for enhanced robustness as reported in several previous studies [3–5, 13, 72, 87–89], where  $K$  is the total number of polynomial coefficients given by:

$$K(r, d) = \frac{1}{d!} \prod_{l=1}^d (r + l), \quad (9)$$

where  $d \in [2, 3]$  is the number of space dimensions. The central stencil  $\mathcal{S}^1$  is given by

$$\mathcal{S}_i^c = \bigcup_{m=0}^{M_c} V_m, \quad (10)$$

where the index  $m$  refers to the local numbering of the elements in the stencil with the element with index 0 being the considered cell  $i$ , and the index  $c$  referring to the stencil number (in case of multiple stencils) where for the central stencil  $c = 1$ . The entire stencil of the considered cell  $i$  is transformed in reference space  $\mathcal{S}_i^c$ , where the  $r^{\text{th}}$  order reconstruction polynomial is an expansion over local polynomial basis functions  $\phi_k(\xi, \eta, \zeta)$  given by:

$$p(\xi, \eta, \zeta) = \sum_{k=0}^K a_k \phi_k(\xi, \eta, \zeta) = \mathbf{U}_0 + \sum_{k=1}^K a_k \phi_k(\xi, \eta, \zeta), \quad (11)$$

where  $\mathbf{U}_0$  corresponds to the vector of conserved variables at the considered cell  $i$ , and  $a_k$  are the degrees of freedom of the polynomial. The degrees of freedom  $a_k$  for the polynomial for each cell  $m$  are obtained by satisfying the condition

that the cell average of the reconstruction polynomial  $p(\xi, \eta, \zeta)$  must be equal to the cell average of the solution  $\mathbf{U}_m$ :

$$\int_{V'_m} p(\xi, \eta, \zeta) d\xi d\eta d\zeta = |V'_m| \mathbf{U}_0 + \sum_{k=1}^K \int_{V'_m} a_k \phi_k d\xi d\eta d\zeta = |V'_m| \mathbf{U}_m, \quad m = 1, \dots, M. \quad (12)$$

It needs to be stressed that since for hexahedral, quadrilateral, prisms and pyramid cells the transformation to reference space into a unit element cell cannot be guaranteed, the basis functions  $\psi_k$  also need to satisfy equation (4). The basis functions employed  $\psi_k$  for all the elements in the stencil are defined as follows:

$$\phi_k(\xi, \eta, \zeta) \equiv \psi_k(\xi, \eta, \zeta) - \frac{1}{|V'_0|} \int_{V'_0} \psi_k d\xi d\eta d\zeta \quad k = 1, 2, \dots, K, \quad (13)$$

and in the present study  $\psi_k$  are Legendre polynomials basis functions. Denoting the integrals of the basis function  $k$  over the cell  $m$  in the stencil, and the vector of right-hand side by  $A_{mk}$  and  $b$  respectively as given by

$$A_{mk} = \int_{V'_m} \phi_k d\xi d\eta d\zeta, \quad b_m = |V'_m|(\mathbf{U}_m - \mathbf{U}_0),$$

the equations for degrees of freedom  $a_k$  can be rewritten in a matrix form as:

$$\sum_{k=1}^K A_{mk} a_k = b_m, \quad m = 1, 2, \dots, M. \quad (14)$$

The resulting linear system is solved by a QR decomposition based on Householder transformation [90] while using a Moore-Penrose pseudo-inverse of  $A_{mk}$  which is only computed once at the beginning of the simulation as detailed in [72].

#### 2.1.1. WENO Scheme

The WENO scheme used in this study, employs a non-linear combination of various reconstruction polynomials from the central stencil and additional directional stencils; each polynomial is weighted according to the smoothness of its solution, and it is based on the approaches of [5, 6, 13, 87]. The polynomials are given as:

$$p_i(\xi, \eta, \zeta)^{\text{weno}} = \sum_{s=1}^{s_i} \omega_s p_s(\xi, \eta, \zeta), \quad (15)$$

where  $s_i$  is the total number of stencils. Substituting back to Eq. (11) for  $p_s(\xi, \eta, \zeta)$ , we obtain the following expression:

$$p_s(\xi, \eta, \zeta) = \sum_{k=0}^K a_k^{(s)} \phi_k(\xi, \eta, \zeta). \quad (16)$$

Using the condition that the sum of all weights is unity, yields:

$$\begin{aligned} p_i(\xi, \eta, \zeta)^{\text{weno}} &= \mathbf{U}_0 + \sum_{k=1}^K \left( \sum_{s=0}^{s_i} \omega_s a_k^{(s)} \right) \phi_k(\xi, \eta, \zeta) \\ &\equiv \mathbf{U}_0 + \sum_{k=1}^K \tilde{a}_k \phi_k(\xi, \eta, \zeta), \end{aligned} \quad (17)$$

where  $\tilde{a}_k$  are the reconstructed degrees of freedom; and the non-linear weight  $\omega_m$  is defined as:

$$\omega_s = \frac{\tilde{\omega}_s}{\sum_{s=1}^{s_t} \tilde{\omega}_s} \quad \text{where} \quad \tilde{\omega}_s = \frac{\lambda_s}{(\epsilon + \mathcal{SI}_s)^b}. \quad (18)$$

The smoothness indicator  $\mathcal{SI}_m$  is given by:

$$\mathcal{SI}_s = \sum_{1 \leq |\beta| \leq r} \int_{V'_0} \left( \mathcal{D}^\beta p_s(\xi, \eta, \zeta) \right)^2 (d\xi, d\eta, d\zeta), \quad (19)$$

where  $\beta$  is a multi-index,  $r$  is the polynomial's order,  $\lambda_m$  is the linear weight. The central stencil is assigned a large linear weight  $\lambda_1$  (to be discussed later), while the remaining directional stencils are assigned a value of  $\lambda_s = 1$  and a value to prevent division by zero of  $\epsilon = 10^{-6}$  is used,  $b = 4$  and  $\mathcal{D}$  is the derivative operator. It must be stressed that the purpose of the linear weights in this WENO type of method is not to recover an optimal polynomial, but rather to give a larger influence to the more symmetric central stencil in regions of smooth gradients as detailed in [5, 6, 13, 87].

The smoothness indicator is a quadratic function of the degrees of freedom ( $a_k^s$ ) and Eq. (19) can be rewritten as:

$$\mathcal{SI}_s = \sum_{k=1}^K a_k^s \left( \sum_{q=1}^K \mathcal{OI}_{kq} a_q^s \right), \quad (20)$$

where the oscillation indication matrix  $\mathcal{OI}_{kq}$  is given by:

$$\mathcal{OI}_{kq} = \sum_{1 \leq |\beta| \leq r} \int_{V'_0} \left( \mathcal{D}^\beta \phi_k(\xi, \eta, \zeta) \right) \left( \mathcal{D}^\beta \phi_q(\xi, \eta, \zeta) \right) (d\xi, d\eta, d\zeta), \quad (21)$$

and can be precomputed and stored at the beginning of the simulation. Additionally the WENO reconstruction can be carried out with respect to the characteristic variables, and the reader is referred to [5, 6, 52] and references therein regarding the implementation. Although the use of characteristic decomposition in reconstruction increases the computational cost of the schemes significantly, previous experiments [91] have demonstrated that in some cases it is necessary in order to avoid spurious oscillations for WENO type of schemes, in particular for cases with strong shocks and contact discontinuities as reported by Harten et al. [92]. For the directional stencils the Type3 algorithm introduced in [72] is employed and in Fig. 2 a typical representation of them can be seen.

### 2.1.2. CWENO Scheme

One of the key ingredients of the CWENO scheme is the combination of an optimal (high-order) polynomial  $p_{opt}$  using the central stencil with lower-order polynomials employing the directional stencils. At the presence of smooth data the optimal polynomial is recovered and therefore the desired-order of accuracy is obtained, whereas at the presence of discontinuous data at least one of the lower-order polynomials arising from the directional stencils could contain smooth data, hence essentially reducing the oscillations in the computed solution. All the polynomials involved are subject to the same requirements as previously set of matching the cell averages of the solution, and all of them are overdetermined linear systems solved with the same constrained least-squares technique. The directional stencils employ the Type3 definition as set in [72]. The computational savings compared to the WENO schemes arise from the reduced size of the directional stencils, and the fact that the directional stencils are contained in the central stencil as it can be seen in Fig. 2. The definition of an optimal polynomial given by:

$$p_{opt}(\xi, \eta, \zeta) = \sum_{s=1}^{s_t} \lambda_s p_s(\xi, \eta, \zeta), \quad (22)$$

where  $s$  is the stencil index, with  $c = 1$  being the central,  $c = 2, 3, \dots$  being the directional,  $s_t$  being the total number of stencils, and  $\lambda_s$  being the linear coefficients for each stencil, whose sum is equal to 1. The  $p_1$  polynomial is



not computed directly, but computed by subtracting the lower-order polynomials from the optimum polynomial as follows:

$$p_1(\xi, \eta, \zeta) = \frac{1}{\lambda_1} \left( p_{opt}(\xi, \eta, \zeta) - \sum_{s=2}^{s_t} \lambda_s p_s(\xi, \eta, \zeta) \right). \quad (23)$$

It needs to be noted that the cost of computing  $p_1$  polynomial is significantly higher than a similar order polynomial, since we need to perform the subtractions of the contributions of the lower-order polynomials. This will be seen later on to have a significant impact on the performance of the schemes. The CWENO reconstruction polynomial is given as a non-linear combination of all the polynomials in the following manner:

$$p(\xi, \eta, \zeta)^{cweno} = \sum_{s=1}^{s_t} \omega_s p_s(\xi, \eta, \zeta), \quad (24)$$

where  $\omega_s$  correspond to the non-linear weights assigned to each polynomial, and in regions with smooth data  $\omega_s \approx \lambda_s$ , hence obtaining the high-order approximation from the central stencil, and in regions of discontinuous solutions the reconstructed solution will be mostly influenced from the lower-order polynomials of the directional stencils. where  $\tilde{a}_k$  are the reconstructed degrees of freedom; and the non-linear weight  $\omega_s$  is defined as:

$$\omega_s = \frac{\tilde{\omega}_s}{\sum_{s=1}^{s_t} \tilde{\omega}_s} \quad \text{where} \quad \tilde{\omega}_s = \frac{\lambda_s}{(\epsilon + \mathcal{ST}_s)^b}. \quad (25)$$

Similarly, to the WENO scheme  $\epsilon = 10^{-6}$  is used and  $b = 4$ . For the present study we employ  $r = 1$  for the directional polynomials resulting in 2nd-order of accuracy, and any arbitrary order of accuracy for the polynomial associated with the central stencil. The smoothness indicators used for CWENO scheme are the same as in the WENO scheme defined previously. It needs to be stressed however that since polynomials of different orders are involved the smoothness indicators are also corresponding to polynomials of different order, and all of them have the same influence on the approximation of the non linear weights. The linear weights are assigned by firstly assigning the non-normalised linear weight for the central stencil  $\lambda_1$  an arbitrary value, and then normalising this as follows:

$$\lambda_1 = 1 - \frac{1}{\lambda_1'}, \quad (26)$$

with the linear weights associated with lower-order polynomials being assigned the same linear weights as follows:

$$\lambda_s = \frac{1 - \lambda_1}{s_t - 1}, \quad (27)$$

where  $s_t$  is the total number of stencils.

### 2.1.3. CWENOZ Scheme

The CWENOZ scheme follows in principle the CWENO scheme, the main difference being the approximation of the non-linear weights. As previously a high-order polynomial is combined with lower-order polynomials arising from the directional stencils also using the Type3 definition as set in [72]. The definition of the optimal polynomial remains the same as before and the CWENOZ reconstruction polynomial is given as a non-linear combination of all the polynomials in the following manner:

$$p(\xi, \eta, \zeta)^{cwenoZ} = \sum_{s=1}^{s_t} \omega_s p_s(\xi, \eta, \zeta), \quad (28)$$

where  $\omega_s$  correspond to the non-linear weights assigned to each polynomial. The characteristic difference is the approximation of the non-linear weights  $\omega_s$ . We mentioned previously that the smoothness indicators for the CWENO scheme arise from polynomials of different orders. The WENOZ component of combining unequal degree poly-

mials as introduced by Borges et al. and Castro et al. [93, 94] is employed in this study, but adapted for unequal polynomials, sized reconstruction stencils and arbitrary elements as has recently been reported by [67, 68]. The non-linear weights are now defined as:

$$\omega_s = \frac{\tilde{\omega}_s}{\sum_{s=1}^{s_I} \tilde{\omega}_s} \quad \text{where} \quad \tilde{\omega}_s = \lambda_s \left( 1 + \frac{\tau}{\epsilon + SI_s} \right). \quad (29)$$

With  $\tau$  being the universal oscillation indicator and taken as the absolute difference between the smoothness indicators as follows:

$$\tau = \left( \frac{\sum_{s=2}^{s_I} |SI_s - SI_1|}{s_I - 1} \right)^b. \quad (30)$$

Similarly to the WENO scheme  $\epsilon = 10^{-6}$  is used and  $b = 4$ . For the present study we employ  $r = 1$  for the directional polynomials resulting in 2nd-order of accuracy, and any arbitrary order of accuracy for the polynomial associated with the central stencil. The procedure for the assignment of the linear weights is similar to the CWENO approach described previously. A more sophisticated way to computing the coefficient  $b$ , could follow the paradigm of Semplice and Visconti [61] and Cravero et al. [95] where a series of comprehensive studies for various parameters, including  $b$  and  $\epsilon$  has been performed for obtaining the optimal convergence rates of the designed schemes for structured meshes. Therefore the expansion of this optimisation to unstructured meshes could further fortify the CWENOZ schemes.

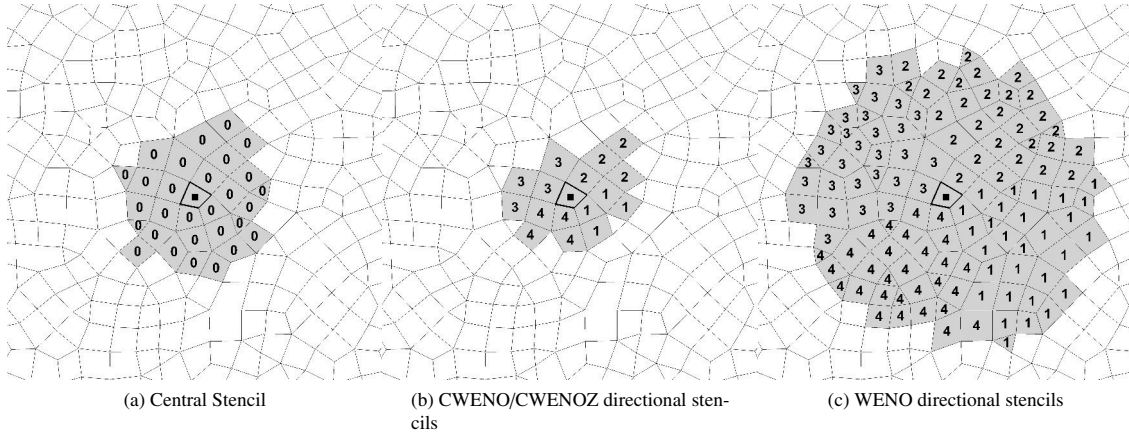


Figure 2: Examples of central and directional stencils for  $\mathcal{P} = 4$  for the WENO, and CWENO/CWENOZ schemes. The considered cell is illustrated by bold lines, the elements of the central stencil indicated by the number 0 and the elements of each of the directional stencils indicated by the numbers 1, 2, 3 and 4. It can be noticed that the CWENO/Z schemes employ significantly smaller directional stencils compared to WENO schemes.

## 2.2. Fluxes approximation & Temporal discretisation

For the inviscid fluxes the approximate HLLC (Harten-Lax-van Leer-Contact) Riemann solver of Toro [96] is employed, unless otherwise stated. For the viscous stress tensor and the heat flux vector in the Navier-Stokes equations the gradients of the discontinuous states for the approximation of the viscous fluxes are averaged by including the penalty terms similar to previous approaches [97–99] in the following manner:

$$\nabla \mathbf{U} = \frac{1}{2} (\nabla \mathbf{U}_L + \nabla \mathbf{U}_R) + \frac{\alpha}{L_{int}} (\mathbf{U}_R - \mathbf{U}_L) \vec{n}, \quad (31)$$

where  $L_{int}$  is the distance between the cell centres of adjacent cells, and  $\alpha = 4/3$  similarly to previous approaches [97, 98]. The gradients are obtained through the unlimited k-exact least square reconstruction utilising the central stencil as previously described in [6, 11], although a weighted least-squares reconstruction could also be employed it was dismissed since from a comprehensive previous study [72] it was documented that it offered a marginal improvement for the solution approximation but no improvement for the gradients approximation. The temporal discretisation employs the 4<sup>th</sup>-order explicit Strong Stability Preserving (SSP) Runge-Kutta method of Spiteri and Ruuth [100] which is stable for  $CFL \approx 1.50$ . The minimum timestep size is defined as follows:

$$\Delta t = \min \left( \frac{1}{d \cdot \left( \frac{s_i}{h_i} + \frac{(\mu+\kappa)}{2 \cdot h_i^2} \right)} \right), \quad (32)$$

where  $d$  refers to the number of dimensions,  $s_i$  is the maximum eigenvalue for the convective flux jacobian as  $s_i = \max(|u_i|, |v_i|, |w_i|) + a_i$ , with  $u, v, w$  being the cartesian velocity components, and  $a_i$  being the speed of sound. Finally, the  $\mu$  and  $\kappa$  correspond to the dynamic viscosity and thermal conductivity coefficient.

All the volume/surface/line integrals are approximated by Gaussian quadrature rule suitable for the order of polynomial employed, while for the initialisation the control volume averages are obtained with a 7th-order Gaussian quadrature to ensure consistent initial conditions across all the schemes. All the schemes developed are implemented in the UCNS3D CFD code [73] which is written in object-oriented Fortran 2003, employing MPI message passing interface (MPI), and the Open Multi-Processing (OpenMP) application programming interface (API), and the reader is referred to [86, 101] for more details on implementation and performance benchmarks. The computational times reported in each case are obtained for the same hardware (and compilation settings) and normalised with respect to a reference setup for the same hardware so that the performance of various algorithms under the same hardware can be appreciated.

### 3. Applications

We present the numerical simulations employed to assess the performance of the CWENO and CWENOZ schemes in terms of robustness, accuracy and computational efficiency for the solution of the Euler and Navier-Stokes equations. Several benchmark test problems have been performed and are listed below:

- 2D Vortex Evolution. This test problem provides an assessment of the accuracy and computational footprint of the methods. The influence of the central stencil linear weight is also investigated.
- 2D Solid Body Rotation. This test problem is ideal for assessing simultaneously the accuracy and non-oscillatory behaviour of the schemes.
- Shu-Osher Problem. This test problem provides an assessment of the non-oscillatory properties of the methods in combination with smooth flow features.
- 2D Riemann Problem. Apart from the non-oscillatory properties of the schemes the impact of the characteristics-based reconstruction is explored for this test case.
- Taylor Green Vortex. A well-established test problem for assessing the ILES capabilities of the schemes. This test is used for establishing the computational cost of all the schemes in a 3D setting.
- Schardin's problem. This test problem is used to assess the performance of a scheme in a flow problem characterised by the presence of strong-gradients regions interacting with vortices and compare with the experimental results.
- Transonic Cylinder  $Re = 3900$ . A test problem involving the interactions of the vortices at the cylinder wake with shocks and therefore being ideal for demonstrating the capability of a numerical scheme to resolve these.

### 3.1. 2D Vortex Evolution

The 2D vortex evolution test problem proposed by Balsara and Shu [102] is used, involving an isentropic vortex propagating at supersonic Mach number at  $45^\circ$  across the domain modelled by the unsteady inviscid Euler equations. The computational domain is given by  $[0, 10] \times [0, 10]$  with periodic boundary conditions applied on all sides. The unperturbed domain has an initial condition  $(\rho, u, v, p) = (1, 1, 1, 1)$ , where temperature and density are defined as  $T = p/\rho$ , and  $S = p/\rho^\gamma$  the adiabatic gas constant  $\gamma = 1.4$  and the vortex perturbations are given by:

$$\delta T = -\frac{(\gamma-1)\epsilon^2}{8\gamma\pi^2}e^{(1-r^2)}, \quad (\delta u, \delta v) = \frac{\epsilon}{2\pi}e^{0.5(1-r^2)}(-(y-5), (x-5)). \quad (33)$$

The vortex strength  $\epsilon = 5$  and adiabatic gas constant  $\gamma = 1.4$ . The  $e_{L^2}$  and the  $e_{L^\infty}$  errors are computed as follows:

$$e_{L^2} = \sqrt{\frac{\sum_i \int_{\Omega_i} (\mathbf{U}_e(x, t_f) - \mathbf{U}_c(x, t_f))^2 dV}{\sum_i |\Omega_i|}}, \quad (34)$$

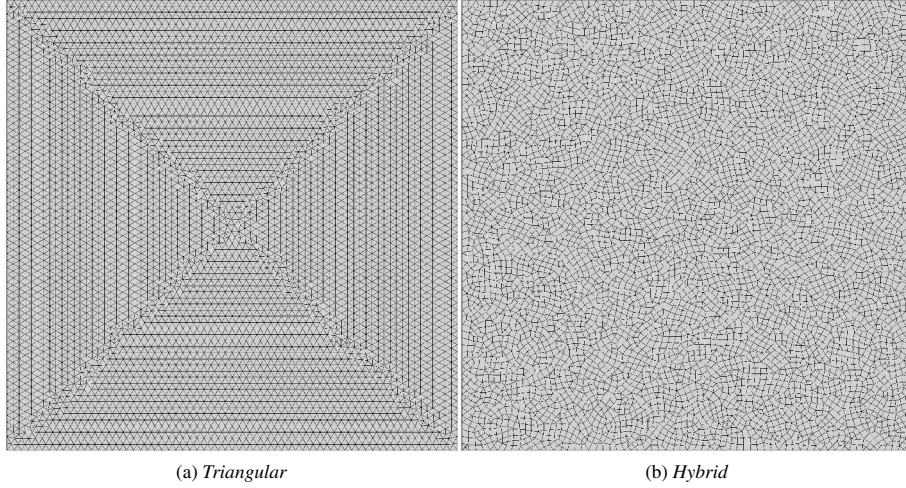


Figure 3: Meshes used for the 2D vortex evolution test problem.

$$e_{L^\infty} = \text{Max} \left| (\mathbf{U}_e(x, t_f) - \mathbf{U}_c(x, t_f)) \right|, \quad (35)$$

where  $\mathbf{U}_c(x, t_f)$  and  $\mathbf{U}_e(x, t_f)$  are the computed and exact solutions at the end of the simulation  $t_f$ . The exact solution  $\mathbf{U}_e(x, t_f)$  being given by the initial condition itself at  $t_0$ . Two types of unstructured meshes a triangular and a hybrid one as shown in Fig. 3 are used for this test problem of 16,32,64 and 128 edges per side resolution, and the simulation is run for a time of  $t_f = 10$ . The purpose of this test problem is to assess the performance of the CWENO and CWENOZ schemes in terms of accuracy, sensitivity to central stencil linear weight, and computational cost against the traditional WENO scheme and unlimited linear schemes. Therefore, schemes ranging from 3rd- to 6th-order of spatial accuracy are employed, and all the WENO variants are assessed using three different values for their central stencil linear weight. To assess the computational cost of the schemes, the time taken for completing each run for the non-linear schemes has been normalised with respect to the time taken by the unlimited linear scheme of equal order of accuracy, at the exact same grid type and resolution and CPU hardware. In this manner a more representative computational footprint of the non-linear schemes is provided. Ideally we would like to determine the smallest value of central stencil linear weight required to achieve the designed order of accuracy, given that WENO, CWENO and CWENOZ will achieve under different conditions. This will be later one used to assess the non-oscillatory properties

of the schemes, so that the central stencil linear weight selected should be suitable to allow the scheme to reach the designed order of spatial accuracy in smooth regions of the flow.

From the obtained results of the unlimited linear schemes as listed in Table 1, it is clear that all the schemes achieve convergence rates close to their theoretical ones. Although for the 5th and 6th-order schemes a timestep size significantly smaller than the one employed with the current time discretisation would have been required to achieve exactly the theoretical ones, but this is beyond the scope of the present study.

Table 1:  $e_{L^\infty}$  and  $e_{L^2}$  error, and convergence rates  $O_{L^\infty}$ ,  $O_{L^2}$  of density for the 2D vortex evolution test problem at the final time of  $t = 10$ , obtained with the linear (unlimited) scheme using various spatial order of accuracy approximations ranging from 3rd- ( $\mathcal{P}2$  Reconstruction Polynomial) to 6th-order ( $\mathcal{P}5$  Reconstruction Polynomial). It can be noticed that all the variants achieve convergence rates close to their theoretical ones.

Order/Number of Edges	Triangular Mesh				Hybrid Mesh			
	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$
$\mathcal{P}2/16$	1.830E-01	0.00	2.266E-02	-	2.828E-01	-	3.620E-02	-
$\mathcal{P}2/32$	3.478E-02	2.40	5.057E-03	2.16	7.661E-02	1.88	1.008E-02	1.84
$\mathcal{P}2/64$	6.089E-03	2.51	9.271E-04	2.45	1.268E-02	2.60	1.878E-03	2.42
$\mathcal{P}2/128$	7.862E-04	2.95	1.257E-04	2.88	1.891E-03	2.74	2.767E-04	2.76
$\mathcal{P}3/16$	3.559E-02	-	5.361E-03	-	1.035E-01	-	1.364E-02	-
$\mathcal{P}3/32$	9.397E-03	1.92	8.548E-04	2.65	2.932E-02	1.82	2.450E-03	2.48
$\mathcal{P}3/64$	3.942E-04	4.58	4.827E-05	4.15	9.213E-04	4.99	1.489E-04	4.04
$\mathcal{P}3/128$	2.737E-05	3.85	3.043E-06	3.99	6.475E-05	3.83	7.436E-06	4.32
$\mathcal{P}4/16$	5.131E-02	-	6.727E-03	-	1.348E-01	-	1.632E-02	-
$\mathcal{P}4/32$	1.557E-02	1.72	1.290E-03	2.38	2.992E-02	2.17	2.702E-03	2.59
$\mathcal{P}4/64$	4.223E-04	5.20	6.881E-05	4.23	1.011E-03	4.89	1.599E-04	4.08
$\mathcal{P}4/128$	1.510E-05	4.81	2.520E-06	4.77	4.188E-05	4.59	6.426E-06	4.64
$\mathcal{P}5/16$	5.532E-02	-	4.804E-03	-	4.396E-02	-	6.456E-03	-
$\mathcal{P}5/32$	2.229E-03	4.63	2.930E-04	4.04	1.357E-02	1.70	1.019E-03	2.66
$\mathcal{P}5/64$	7.700E-05	4.86	8.411E-06	5.12	2.307E-04	5.88	2.566E-05	5.31
$\mathcal{P}5/128$	1.672E-06	5.53	4.616E-07	4.19	3.946E-06	5.87	6.382E-07	5.33

From the WENO schemes results listed in Table 2, all the variants reach convergence rates close to their theoretical ones and similar errors to the unlimited linear schemes. The obtained convergence rates are not particularly sensitive to the choice of the central stencil linear weight as also previously documented in [13, 14]. The computational cost of the WENO schemes is approximately 3.4 times higher than the equivalent unlimited scheme. From this study a linear weight of  $\lambda_1 = 10^3$  will be used hereafter for WENO scheme.

From the CWENO schemes results listed in Table 3, it can be noticed that the convergence rates are particularly sensitive to the choice of the central stencil linear weight. Firstly the linear weight  $\lambda_1 = 10^3$  is clearly unsuitable for all discretisation orders and mesh resolutions. The linear weight  $\lambda_1 = 10^7$  is providing a better performance compared to the smaller one, but as the spatial order of accuracy of the scheme increases the performance deteriorated. On the other hand, with an excessively large central stencil linear weight of  $\lambda_1 = 10^{15}$  the obtained performance is close to the equivalent unlimited linear scheme. The unscaled smoothness indicators arising from polynomials of different orders of accuracy prevent the CWENO schemes from realising their full potential (especially for higher-order polynomials), unless an excessive large central stencil linear weight is used. On the positive side the computational cost of the CWENO schemes is approximately 2.4 times higher than the equivalent unlimited scheme, therefore establishing them as preferred alternative to WENO schemes in terms of cost. From this study a linear weight of  $\lambda_1 = 10^{15}$  will be the selected one for the CWENO scheme.

The CWENOZ schemes achieve convergence rates close to the theoretical ones as shown in Table 4, while not being sensitive to the choice of the central stencil linear weight. It needs to be highlighted that this desirable insensi-

Table 2:  $e_{L^\infty}$  and  $e_{L^2}$  error, convergence rates  $O_{L^\infty}$ ,  $O_{L^2}$  of density, and normalised CPU cost (normalised with respect to the CPU time taken for the equivalent unlimited scheme at the same grid resolution) for the 2D vortex evolution test problem at the final time of  $t = 10$  obtained with the WENO schemes using various spatial order of accuracy approximations ranging from 3rd-( $\mathcal{P}2$  Reconstruction Polynomial) to 6th-order ( $\mathcal{P}5$  Reconstruction Polynomial), and linear weight  $\lambda_1$  for the central stencil. It can be noticed that all the variants achieve convergence rates close to their theoretical ones, the dependence of the convergence rates to the central stencil linear weight is limited, while being approximately 3.4 times more expensive than the equivalent unlimited scheme.

Order/Number of Edges	$\lambda_1 = 10^3$				$\lambda_1 = 10^6$				$\lambda_1 = 10^9$				Cost
Triangular Mesh	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	CPU
$\mathcal{P}2/16$	1.934E-01	-	2.449E-02	-	1.833E-01	-	2.272E-02	-	1.830E-01	-	2.266E-02	-	3.326
$\mathcal{P}2/32$	3.430E-02	2.50	4.881E-03	2.33	3.478E-02	2.40	5.056E-03	2.17	3.478E-02	2.40	5.057E-03	2.16	2.947
$\mathcal{P}2/64$	6.059E-03	2.50	9.202E-04	2.41	6.089E-03	2.51	9.271E-04	2.45	6.089E-03	2.51	9.271E-04	2.45	3.773
$\mathcal{P}2/128$	7.818E-04	2.95	1.249E-04	2.88	7.862E-04	2.95	1.257E-04	2.88	7.862E-04	2.95	1.257E-04	2.88	3.412
$\mathcal{P}3/16$	4.107E-02	-	6.814E-03	-	3.633E-02	-	5.422E-03	-	3.568E-02	-	5.364E-03	-	3.033
$\mathcal{P}3/32$	9.258E-03	2.15	8.254E-04	3.05	9.398E-03	1.95	8.547E-04	2.67	9.397E-03	1.92	8.548E-04	2.65	3.195
$\mathcal{P}3/64$	3.900E-04	4.57	4.761E-05	4.12	3.942E-04	4.58	4.827E-05	4.15	3.942E-04	4.58	4.827E-05	4.15	3.051
$\mathcal{P}3/128$	2.708E-05	3.85	3.013E-06	3.98	2.737E-05	3.85	3.043E-06	3.99	2.737E-05	3.85	3.043E-06	3.99	2.999
$\mathcal{P}4/16$	6.070E-02	-	8.459E-03	-	5.129E-02	-	6.720E-03	-	5.132E-02	-	6.727E-03	-	3.224
$\mathcal{P}4/32$	1.550E-02	1.97	1.285E-03	2.72	1.557E-02	1.72	1.290E-03	2.38	1.557E-02	1.72	1.290E-03	2.38	3.158
$\mathcal{P}4/64$	4.222E-04	5.20	6.876E-05	4.22	4.223E-04	5.20	6.881E-05	4.23	4.223E-04	5.20	6.881E-05	4.23	3.483
$\mathcal{P}4/128$	1.513E-05	4.80	2.524E-06	4.77	1.510E-05	4.81	2.520E-06	4.77	1.510E-05	4.81	2.520E-06	4.77	3.183
$\mathcal{P}5/16$	3.887E-02	-	4.751E-03	-	5.537E-02	-	4.846E-03	-	5.533E-02	-	4.803E-03	-	3.299
$\mathcal{P}5/32$	2.238E-03	4.12	2.959E-04	4.01	2.229E-03	4.63	2.930E-04	4.05	2.229E-03	4.63	2.930E-04	4.04	3.249
$\mathcal{P}5/64$	7.729E-05	4.86	8.494E-06	5.12	7.700E-05	4.86	8.411E-06	5.12	7.700E-05	4.86	8.411E-06	5.12	3.541
$\mathcal{P}5/128$	1.686E-06	5.52	4.620E-07	4.20	1.672E-06	5.53	4.616E-07	4.19	1.672E-06	5.53	4.616E-07	4.19	3.663
Hybrid Mesh	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	CPU
$\mathcal{P}2/16$	3.177E-01	-	4.259E-02	-	2.914E-01	-	3.767E-02	-	2.839E-01	-	3.641E-02	-	3.364
$\mathcal{P}2/32$	8.134E-02	1.97	9.919E-03	2.10	7.661E-02	1.93	1.005E-02	1.91	7.661E-02	1.89	1.008E-02	1.85	3.518
$\mathcal{P}2/64$	1.244E-02	2.71	1.809E-03	2.46	1.267E-02	2.60	1.878E-03	2.42	1.268E-02	2.60	1.878E-03	2.42	3.482
$\mathcal{P}2/128$	1.871E-03	2.73	2.730E-04	2.73	1.891E-03	2.74	2.767E-04	2.76	1.891E-03	2.74	2.767E-04	2.76	3.005
$\mathcal{P}3/16$	1.638E-01	-	2.468E-02	-	1.151E-01	-	1.604E-02	-	1.044E-01	-	1.386E-02	-	3.187
$\mathcal{P}3/32$	2.743E-02	2.58	2.226E-03	3.47	2.923E-02	1.98	2.429E-03	2.72	2.931E-02	1.83	2.449E-03	2.50	3.647
$\mathcal{P}3/64$	8.979E-04	4.93	1.441E-04	3.95	9.212E-04	4.99	1.489E-04	4.03	9.213E-04	4.99	1.489E-04	4.04	3.352
$\mathcal{P}3/128$	6.263E-05	3.84	7.166E-06	4.33	6.474E-05	3.83	7.436E-06	4.32	6.475E-05	3.83	7.436E-06	4.32	3.314
$\mathcal{P}4/16$	1.747E-01	-	2.308E-02	-	1.415E-01	-	1.763E-02	-	1.358E-01	-	1.647E-02	-	4.422
$\mathcal{P}4/32$	2.939E-02	2.57	2.617E-03	3.14	2.990E-02	2.24	2.697E-03	2.71	2.992E-02	2.18	2.702E-03	2.61	3.938
$\mathcal{P}4/64$	1.015E-03	4.86	1.599E-04	4.03	1.011E-03	4.89	1.599E-04	4.08	1.011E-03	4.89	1.599E-04	4.08	3.880
$\mathcal{P}4/128$	4.239E-05	4.58	6.497E-06	4.62	4.188E-05	4.59	6.426E-06	4.64	4.188E-05	4.59	6.426E-06	4.64	3.618
$\mathcal{P}5/16$	1.003E-01	-	1.397E-02	-	5.540E-02	-	7.847E-03	-	4.449E-02	-	6.580E-03	-	3.598
$\mathcal{P}5/32$	1.270E-02	2.98	9.734E-04	3.84	1.357E-02	2.03	1.016E-03	2.95	1.357E-02	1.71	1.018E-03	2.69	3.878
$\mathcal{P}5/64$	2.161E-04	5.88	2.635E-05	5.21	2.307E-04	5.88	2.566E-05	5.31	2.307E-04	5.88	2.566E-05	5.31	4.180
$\mathcal{P}5/128$	4.177E-06	5.69	6.582E-07	5.32	3.946E-06	5.87	6.382E-07	5.33	3.946E-06	5.87	6.382E-07	5.33	4.199

tivity to the choice of the central stencil linear weight is the result of the scaling of the smoothness indicators arising from polynomials of different orders of accuracy. The computational cost of the CWENOZ schemes is similar to the CWENO schemes and approximately 2.4 times higher than the equivalent unlimited scheme. From this study a central stencil linear weight of  $\lambda_1 = 0.9$  will be the selected one for CWENOZ.

### 3.2. 2D Solid Body Rotation

The solid body rotation test of Leveque [103] is employed to investigate the performance of the WENO, CWENO and CWENOZ schemes in terms of their non-oscillatory properties on the correct and accurate revolution of a body. The continuity equation is considered as follows:

$$\frac{\partial U}{\partial t} + \nabla \cdot (\mathbf{v}U) = 0, \quad (36)$$



Table 3:  $e_{L^\infty}$  and  $e_{L^2}$  error, convergence rates  $O_{L^\infty}$ ,  $O_{L^2}$  of density, and normalised CPU cost (normalised with respect to the CPU time taken for the equivalent unlimited scheme at the same grid resolution) for the 2D vortex evolution test problem at the final time of  $t = 10$  obtained with the CWENO schemes using various spatial order of accuracy approximations ranging from 3rd-( $\mathcal{P}2$  Reconstruction Polynomial) to 6th-order ( $\mathcal{P}5$  Reconstruction Polynomial), and linear weight  $\lambda_1$  for the central stencil. It can be noticed the scheme achieves convergence rates close to their theoretical one for all grid resolutions only when the largest value for the central stencil linear weight is used and is approximately 2.4 times more expensive than the equivalent unlimited scheme.

Order/Number of Edges	$\lambda_1 = 10^3$				$\lambda_1 = 10^7$				$\lambda_1 = 10^{15}$				Cost
Triangular Mesh	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	CPU
$\mathcal{P}2/16$	3.166E-01	-	4.117E-02	-	1.963E-01	-	2.424E-02	-	1.832E-01	-	2.268E-02	-	2.743
$\mathcal{P}2/32$	1.338E-01	1.24	1.467E-02	1.49	4.059E-02	2.27	5.591E-03	2.12	3.477E-02	2.40	5.059E-03	2.16	1.797
$\mathcal{P}2/64$	3.296E-02	2.02	3.256E-03	2.17	7.734E-03	2.39	1.036E-03	2.43	6.088E-03	2.51	9.269E-04	2.45	2.331
$\mathcal{P}2/128$	7.643E-03	2.11	4.475E-04	2.86	7.785E-04	3.31	1.266E-04	3.03	7.862E-04	2.95	1.257E-04	2.88	2.732
$\mathcal{P}3/16$	3.321E-01	-	4.292E-02	-	1.590E-01	-	1.759E-02	-	3.542E-02	-	5.385E-03	-	2.596
$\mathcal{P}3/32$	1.486E-01	1.16	1.538E-02	1.48	3.492E-02	2.19	3.946E-03	2.16	9.499E-03	1.90	8.616E-04	2.64	2.822
$\mathcal{P}3/64$	5.092E-02	1.55	5.508E-03	1.48	1.316E-02	1.41	7.735E-04	2.35	3.921E-04	4.60	4.784E-05	4.17	2.875
$\mathcal{P}3/128$	1.684E-02	1.60	1.302E-03	2.08	1.321E-03	3.32	3.898E-05	4.31	2.737E-05	3.84	3.043E-06	3.97	2.621
$\mathcal{P}4/16$	3.389E-01	-	4.416E-02	-	2.324E-01	-	2.833E-02	-	5.209E-02	-	6.778E-03	-	2.561
$\mathcal{P}4/32$	1.445E-01	1.23	1.640E-02	1.43	8.291E-02	1.49	8.842E-03	1.68	1.578E-02	1.72	1.306E-03	2.38	2.668
$\mathcal{P}4/64$	5.196E-02	1.48	5.381E-03	1.61	1.984E-02	2.06	1.387E-03	2.67	4.206E-04	5.23	6.855E-05	4.25	2.624
$\mathcal{P}4/128$	1.949E-02	1.41	1.405E-03	1.94	8.909E-04	4.48	3.534E-05	5.29	1.510E-05	4.80	2.520E-06	4.77	2.518
$\mathcal{P}5/16$	3.472E-01	-	4.551E-02	-	2.760E-01	-	3.627E-02	-	5.605E-02	-	4.949E-03	-	2.939
$\mathcal{P}5/32$	1.590E-01	1.13	1.861E-02	1.29	1.193E-01	1.21	1.252E-02	1.53	2.413E-03	4.54	3.107E-04	3.99	2.325
$\mathcal{P}5/64$	6.346E-02	1.32	6.274E-03	1.57	4.412E-02	1.44	3.949E-03	1.66	7.198E-05	5.07	8.041E-06	5.27	2.404
$\mathcal{P}5/128$	2.502E-02	1.34	2.003E-03	1.65	9.912E-03	2.15	5.227E-04	2.92	1.672E-06	5.43	4.616E-07	4.12	2.354
Hybrid Mesh	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	CPU
$\mathcal{P}2/16$	3.775E-01	-	5.253E-02	-	3.101E-01	-	4.077E-02	-	2.825E-01	-	3.615E-02	-	2.859
$\mathcal{P}2/32$	2.260E-01	0.74	2.694E-02	0.96	9.416E-02	1.72	1.102E-02	1.89	7.682E-02	1.88	1.009E-02	1.84	3.880
$\mathcal{P}2/64$	6.290E-02	1.85	6.829E-03	1.98	1.809E-02	2.38	2.164E-03	2.35	1.271E-02	2.60	1.877E-03	2.43	3.303
$\mathcal{P}2/128$	1.867E-02	1.75	1.345E-03	2.34	2.207E-03	3.04	2.873E-04	2.91	1.891E-03	2.75	2.767E-04	2.76	2.764
$\mathcal{P}3/16$	3.837E-01	-	5.367E-02	-	2.748E-01	-	3.736E-02	-	1.012E-01	-	1.332E-02	-	2.676
$\mathcal{P}3/32$	2.415E-01	0.67	2.857E-02	0.91	9.248E-02	1.57	1.103E-02	1.76	2.924E-02	1.79	2.426E-03	2.46	3.020
$\mathcal{P}3/64$	7.839E-02	1.62	8.142E-03	1.81	3.173E-02	1.54	2.198E-03	2.33	9.806E-04	4.90	1.499E-04	4.02	3.057
$\mathcal{P}3/128$	3.065E-02	1.35	2.884E-03	1.50	3.311E-03	3.26	2.012E-04	3.45	6.475E-05	3.92	7.436E-06	4.33	2.587
$\mathcal{P}4/16$	3.880E-01	-	5.443E-02	-	3.302E-01	-	4.446E-02	-	1.291E-01	-	1.552E-02	-	2.769
$\mathcal{P}4/32$	2.470E-01	0.65	3.099E-02	0.81	1.494E-01	1.14	1.584E-02	1.49	3.056E-02	2.08	2.670E-03	2.54	2.817
$\mathcal{P}4/64$	8.488E-02	1.54	8.766E-03	1.82	4.024E-02	1.89	3.605E-03	2.14	9.727E-04	4.97	1.604E-04	4.06	2.763
$\mathcal{P}4/128$	3.417E-02	1.31	3.224E-03	1.44	3.576E-03	3.49	2.926E-04	3.62	4.189E-05	4.54	6.424E-06	4.64	2.445
$\mathcal{P}5/16$	3.915E-01	-	5.493E-02	-	3.600E-01	-	4.940E-02	-	3.820E-02	-	7.712E-03	-	2.313
$\mathcal{P}5/32$	2.639E-01	0.57	3.389E-02	0.70	2.020E-01	0.83	2.207E-02	1.16	1.337E-02	1.51	1.066E-03	2.86	2.490
$\mathcal{P}5/64$	9.875E-02	1.42	9.921E-03	1.77	6.608E-02	1.61	6.921E-03	1.67	8.712E-04	3.94	4.394E-05	4.60	2.575
$\mathcal{P}5/128$	3.759E-02	1.39	3.959E-03	1.33	1.869E-02	1.82	1.625E-03	2.09	3.966E-06	7.78	6.381E-07	6.11	2.339

on a computational domain given by  $[0, 1]^2$  with periodic boundary conditions applied on all sides. The velocity  $\mathbf{v}$  describing the revolution of the profile around the centre of the domain  $(0.5, 0.5)$  is given by:

$$\mathbf{v}(x, y) = (0.5 - y, x - 0.5). \quad (37)$$

In this particular test three bodies are considered namely a smooth centered at  $(x_0 = 0.25, y_0 = 0.5)$ , a sharp cone centered at  $(x_0 = 0.5, y_0 = 0.25)$  and a slotted cylinder centered at  $(x_0 = 0.5, y_0 = 0.75)$  described by the following functions respectively:

$$\mathbf{f}(x, y) = \frac{1 + \cos(\pi r(x, y))}{4}, \quad (38)$$

$$\mathbf{f}(x, y) = 1 - r(x, y), \quad (39)$$

and

$$\mathbf{f}(x, y) = \begin{cases} 1, & \text{if } |x - x_0| \geq 0.025 \text{ or } y \geq 0.85 \\ 0, & \text{if otherwise.} \end{cases} \quad (40)$$

Table 4:  $e_{L^\infty}$  and  $e_{L^2}$  error, convergence rates  $O_{L^\infty}$ ,  $O_{L^2}$  of density, and normalised CPU cost (normalised with respect to the CPU time taken for the equivalent unlimited scheme at the same grid resolution) for the 2D vortex evolution test problem at the final time of  $t = 10$  obtained with the CWENOZ schemes using various spatial order of accuracy approximations ranging from 3rd- ( $\mathcal{P}2$  Reconstruction Polynomial) to 6th-order ( $\mathcal{P}5$  Reconstruction Polynomial), and linear weight  $\lambda_1$  for the central stencil. It can be noticed the scheme achieves convergence rates close to their theoretical while not being significantly influenced by the central stencil linear weight value, and is approximately 2.4 times more expensive than the equivalent unlimited scheme.

Order/Number of Edges	$\lambda_1 = 0.9$				$\lambda_1 = 0.999$				$\lambda_1 = 0.999999$				Cost
Triangular Mesh	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	CPU
$\mathcal{P}2/16$	1.779E-01	-	2.174E-02	-	1.829E-01	-	2.264E-02	-	1.830E-01	-	2.266E-02	-	2.676
$\mathcal{P}2/32$	3.475E-02	2.36	5.055E-03	2.10	3.478E-02	2.40	5.057E-03	2.16	3.478E-02	2.40	5.057E-03	2.16	1.791
$\mathcal{P}2/64$	6.089E-03	2.51	9.271E-04	2.45	6.089E-03	2.51	9.271E-04	2.45	6.089E-03	2.51	9.271E-04	2.45	2.192
$\mathcal{P}2/128$	7.862E-04	2.95	1.257E-04	2.88	7.862E-04	2.95	1.257E-04	2.88	7.862E-04	2.95	1.257E-04	2.88	2.624
$\mathcal{P}3/16$	2.018E-01	-	1.868E-02	-	3.507E-02	-	5.272E-03	-	3.559E-02	-	5.361E-03	-	2.567
$\mathcal{P}3/32$	9.936E-03	4.34	8.703E-04	4.42	9.400E-03	1.90	8.549E-04	2.62	9.397E-03	1.92	8.548E-04	2.65	2.733
$\mathcal{P}3/64$	3.942E-04	4.66	4.827E-05	4.17	3.942E-04	4.58	4.827E-05	4.15	3.942E-04	4.58	4.827E-05	4.15	2.558
$\mathcal{P}3/128$	2.737E-05	3.85	3.043E-06	3.99	2.737E-05	3.85	3.043E-06	3.99	2.737E-05	3.85	3.043E-06	3.99	2.530
$\mathcal{P}4/16$	1.449E-01	-	1.760E-02	-	4.806E-02	-	5.915E-03	-	5.124E-02	-	6.720E-03	-	2.471
$\mathcal{P}4/32$	1.716E-02	3.08	1.338E-03	3.72	1.559E-02	1.62	1.290E-03	2.20	1.557E-02	1.72	1.290E-03	2.38	2.428
$\mathcal{P}4/64$	4.223E-04	5.34	6.881E-05	4.28	4.223E-04	5.21	6.881E-05	4.23	4.223E-04	5.20	6.881E-05	4.23	2.440
$\mathcal{P}4/128$	1.510E-05	4.81	2.520E-06	4.77	1.510E-05	4.81	2.520E-06	4.77	1.510E-05	4.81	2.520E-06	4.77	2.373
$\mathcal{P}5/16$	9.085E-02	-	1.943E-02	-	8.923E-02	-	7.104E-03	-	5.540E-02	-	4.804E-03	-	2.451
$\mathcal{P}5/32$	3.099E-03	4.87	3.143E-04	5.95	2.233E-03	5.32	2.925E-04	4.60	2.229E-03	4.64	2.930E-04	4.04	2.254
$\mathcal{P}5/64$	7.699E-05	5.33	8.410E-06	5.22	7.700E-05	4.86	8.411E-06	5.12	7.700E-05	4.86	8.411E-06	5.12	2.308
$\mathcal{P}5/128$	1.672E-06	5.53	4.616E-07	4.19	1.672E-06	5.53	4.616E-07	4.19	1.672E-06	5.53	4.616E-07	4.19	2.306
Hybrid Mesh	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	$e_{L^\infty}$	$O_{L^\infty}$	$e_{L^2}$	$O_{L^2}$	CPU
$\mathcal{P}2/16$	2.720E-01	-	3.436E-02	-	2.823E-01	-	3.611E-02	-	2.828E-01	-	3.620E-02	-	3.456
$\mathcal{P}2/32$	7.631E-02	1.83	9.980E-03	1.78	7.661E-02	1.88	1.008E-02	1.84	7.661E-02	1.88	1.008E-02	1.84	3.244
$\mathcal{P}2/64$	1.268E-02	2.59	1.878E-03	2.41	1.268E-02	2.60	1.878E-03	2.42	1.268E-02	2.60	1.878E-03	2.42	3.184
$\mathcal{P}2/128$	1.891E-03	2.75	2.767E-04	2.76	1.891E-03	2.74	2.767E-04	2.76	1.891E-03	2.74	2.767E-04	2.76	2.737
$\mathcal{P}3/16$	9.570E-02	-	1.697E-02	-	9.008E-02	-	1.186E-02	-	1.035E-01	-	1.363E-02	-	3.323
$\mathcal{P}3/32$	5.191E-02	0.88	3.871E-03	2.13	2.968E-02	1.60	2.469E-03	2.26	2.932E-02	1.82	2.450E-03	2.48	2.977
$\mathcal{P}3/64$	9.206E-04	5.82	1.488E-04	4.70	9.213E-04	5.01	1.489E-04	4.05	9.213E-04	4.99	1.489E-04	4.04	2.943
$\mathcal{P}3/128$	6.475E-05	3.83	7.436E-06	4.32	6.475E-05	3.83	7.436E-06	4.32	6.475E-05	3.83	7.436E-06	4.32	2.577
$\mathcal{P}4/16$	1.353E-01	-	2.077E-02	-	1.136E-01	-	1.397E-02	-	1.347E-01	-	1.630E-02	-	2.873
$\mathcal{P}4/32$	7.738E-02	0.81	5.394E-03	1.94	3.167E-02	1.84	2.757E-03	2.34	2.992E-02	2.17	2.703E-03	2.59	2.761
$\mathcal{P}4/64$	1.010E-03	6.26	1.598E-04	5.08	1.011E-03	4.97	1.599E-04	4.11	1.011E-03	4.89	1.599E-04	4.08	2.752
$\mathcal{P}4/128$	4.188E-05	4.59	6.426E-06	4.64	4.188E-05	4.59	6.426E-06	4.64	4.188E-05	4.59	6.426E-06	4.64	2.437
$\mathcal{P}5/16$	1.872E-01	-	2.655E-02	-	5.940E-02	-	1.192E-02	-	4.354E-02	-	6.389E-03	-	2.343
$\mathcal{P}5/32$	9.904E-02	0.92	7.801E-03	1.77	1.471E-02	2.01	1.084E-03	3.46	1.358E-02	1.68	1.019E-03	2.65	2.476
$\mathcal{P}5/64$	2.300E-04	8.75	2.548E-05	8.26	2.307E-04	5.99	2.566E-05	5.40	2.307E-04	5.88	2.566E-05	5.31	2.526
$\mathcal{P}5/128$	3.946E-06	5.86	6.382E-07	5.32	3.946E-06	5.87	6.382E-07	5.33	3.946E-06	5.87	6.382E-07	5.33	2.336

Where the normalised distance from the centres  $(x_0, y_0)$  is given by:

$$r(x, y) = (1/r_0) \sqrt{(x - x_0)^2 + (y - y_0)^2}, \quad (41)$$

with  $r_0 = 0.15$ . The rest of the domain the solution is initialised with zero, and after one full revolution  $t_f = 2\pi$  the exact solution coincides with the initial solution. A triangular unstructured mesh is used as shown in Fig. 3 with 64 edges per side of the computational domain. The WENO, CWENO and CWENOZ schemes ranging from 3rd- to 6th-order of spatial accuracy are employed, with their selected weights from the previous study. The obtained results following one revolution are illustrated in Fig. 4.

From the obtained results it can be noticed that all schemes capture the correct shapes but oscillations in the solutions are present. The CWENO scheme results exhibit smaller oscillations compared to the other schemes while being accurate at smooth extrema. Overall the WENO and CWENOZ schemes are susceptible to larger oscillations in particular at the base of the slotted cylinder, while the CWENO being more dissipative at the top of the slotted cylinder.



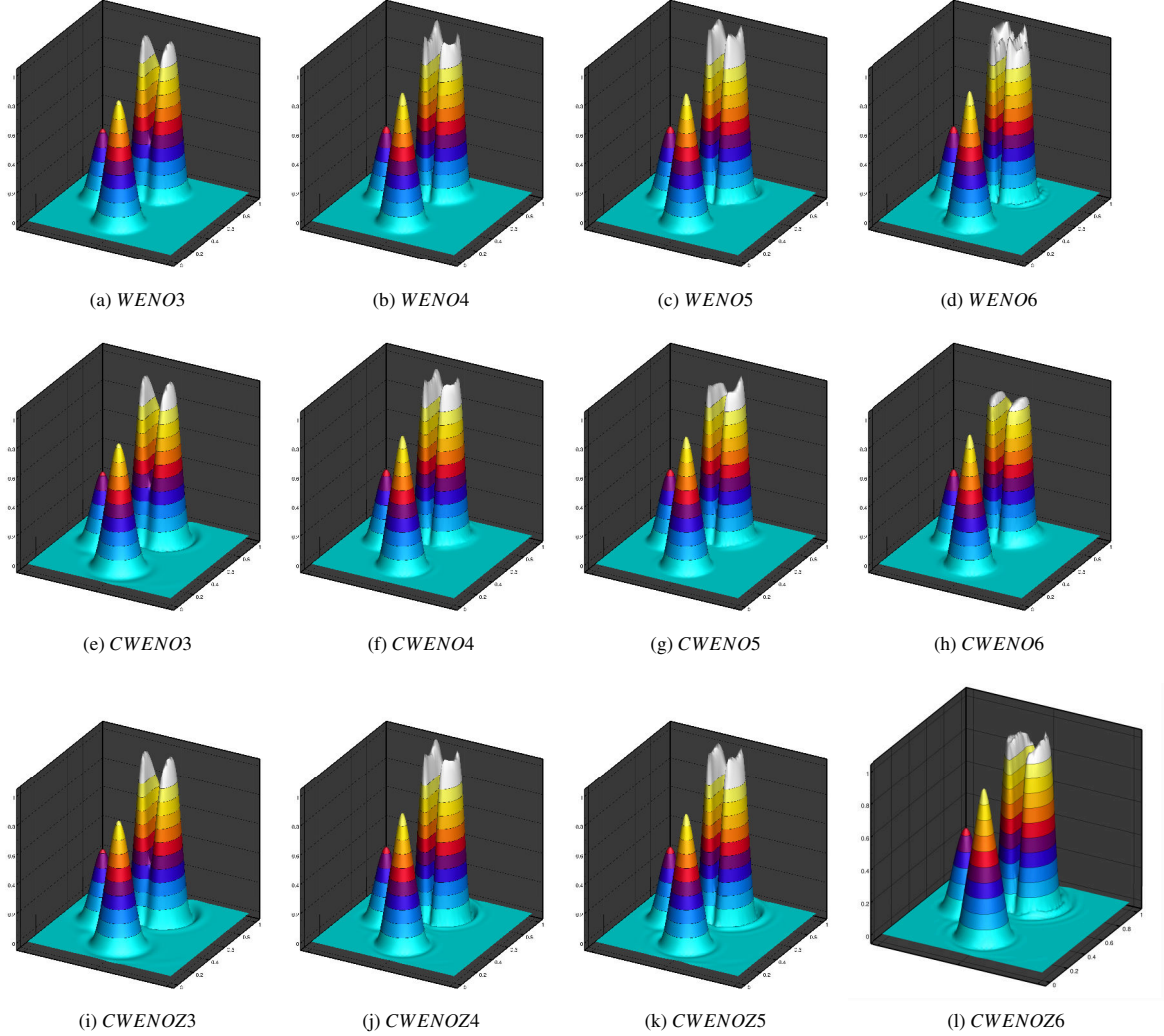


Figure 4: Computed results after one revolution ( $t = 2\pi$ ) of the solid body rotation test problem, with all the schemes. Three-dimensional elevations of the solution coloured by ten uniformly distributed isolines from 0.1 to 1.0. It can be noticed that all schemes capture the correct shapes but oscillations in the solutions are present. The CWENO scheme results exhibit smaller oscillations compared to the other schemes, in particular at the base of the slotted cylinder.

### 3.3. Shu-Osher Problem

The well established Shu-Osher [104] test problem is employed, which involves the interaction between a shock wave and an entropy wave. The computational domain is  $[-4.5, 4.5] \times [0, 1]$ , with periodic boundary conditions in y-axis, a supersonic inflow and outflow on the left and right side of the domain respectively. The initial profile consists of a shock wave  $(\rho, u, v, p) = (3.857143, 2.629369, 0, 10.333333)$  on the left when  $x < -4$  and an entropy wave in the rest of the domain  $(\rho, u, v, p) = (1 + 0.2\sin(5x), 0, 0, 1)$ .

A 2D triangular mesh is utilised with a resolution of each edge being  $h = 0.025$  with approximately 33,886 elements. The reference solution is computed with a one-dimensional solver of the Euler equations using 10,000 grid points and employing a 5th order WENO scheme. The calculation is run until  $t = 1.8$ .

From the density distribution plots shown in Fig. 5 it can be noticed that all the schemes achieve a good agreement with the reference solution, however the WENO 4th-order scheme exhibits a small overprediction of the peaks and valleys at the shock-wave entropy interaction region as also noticed at a smaller extent for CWENO 5th-order scheme. It needs to be highlighted that switching to a higher order CWENO or CWENOZ schemes is considerably cheaper in terms of computational resources compared to the WENO schemes as presented in Table 5. Overall the CWENO and CWENOZ provide a small positive difference compared to the traditional WENO schemes.

Table 5: Normalised computational times (with respect to WENO3) for the Shu-Osher [104] shock tube test problem.

Polynomial Order	WENO	CWENO	CWENOZ
$\mathcal{P}2$	1.00	1.00	0.93
$\mathcal{P}3$	1.44	1.20	1.21
$\mathcal{P}4$	2.25	1.52	1.60

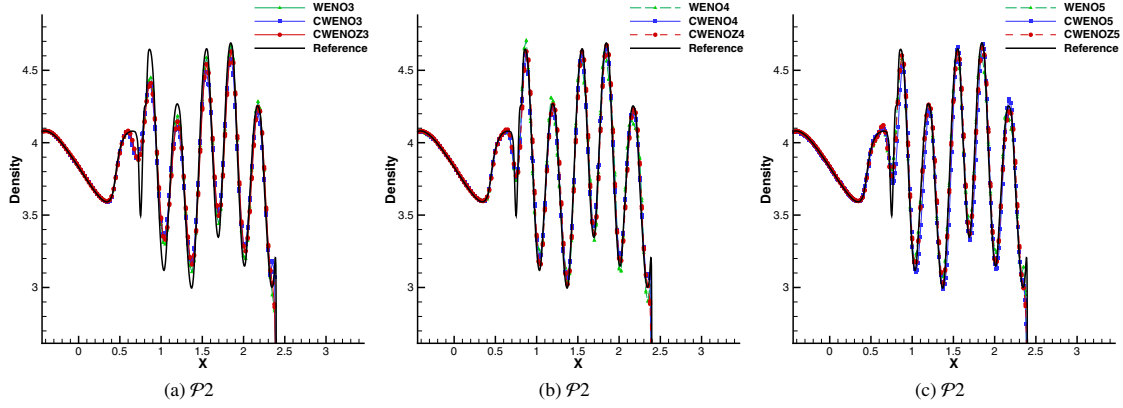


Figure 5: Density distribution for the Shu-Osher [104] shock tube test problem at the final time  $t = 1.8$  using various schemes and comparison with the reference solution obtained from the one-dimensional Euler equations on 10,000 grid points using a WENO-5th order scheme. It can be noticed that the CWENO and CWENOZ are slightly less oscillatory compared to the WENO scheme.

### 3.4. 2D Riemann Problem

The 2D Riemann problem introduced by Schulz et al. [105] is used, where the computational domain is  $[-0.5, 0.5] \times [-0.5, 0.5]$  and the initial condition is given by four different states assigned to each of the quadrants of the domain as shown in Fig. 6. The computational domain has been discretised by a uniform quadrilateral mesh of  $400 \times 400$ . The simulation has been performed until  $t = 0.5$ . The variants of the numerical schemes employed are the 4th-order and 6th-order with the reconstruction carried out in conservative variables, and a 6th-order with the reconstruction carried out in characteristic variables.

From the obtained results in Fig. 7 it can be noticed that all the schemes predict the correct solution pattern, but the WENO implementation provides significantly more oscillatory results compared to the other implementations. This behaviour is expected since all the reconstruction polynomials are of the same order of accuracy, and when a directional stencil has the largest non-linear weight the resulting solution is expected to produce that compared to the CWENO schemes where their directional stencils are lower-order. The WENO scheme results are also the ones with slightly increased number of small-scale structures, and asymmetry of the resulting solution compared to the other schemes. For all the schemes when the reconstruction is carried out in characteristic variables there is a

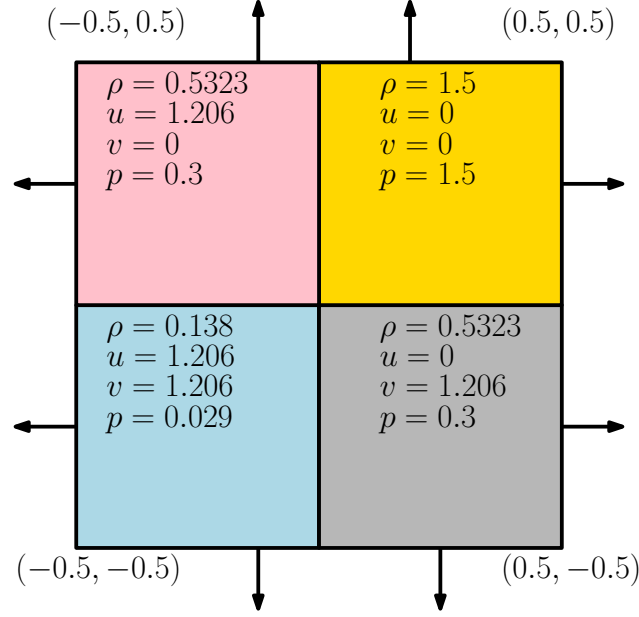


Figure 6: Initial conditions for the multidimensional Riemann problem.

marginal improvement in terms of the roll-ups of the Kelvin-Helmholtz instabilities and the number of structures in the mushroom stem.

### 3.5. ILES of Taylor Green Vortex $Re = 1600$

The ILES of the 3D viscous Taylor-Green vortex test problem at  $Re = 1600$  is employed, for assessing the performance of all the schemes. It is a widely used test problem for the validation of numerical methods, and in particular at relative coarse-“under-resolved” meshes within the LES context [6, 27, 29, 85, 106–110] due to the pronounced dissipation and dispersion characteristics of non-linear methods. The computational domain is defined as  $\Omega = [0, 2\pi]^3$  with periodic boundary conditions. This formulation of the Taylor-Green vortex problem is initialized with the following velocity, density and pressure fields:

$$u(x, y, z, 0) = \sin(kx) \cos(ky) \cos(kz), \quad (42)$$

$$v(x, y, z, 0) = -\cos(kx) \sin(ky) \cos(kz), \quad (43)$$

$$w(x, y, z, 0) = 0, \quad (44)$$

$$\rho(x, y, z, 0) = 1, \quad (45)$$

$$p(x, y, z, 0) = 100 + \frac{P}{16} [\cos(2z) + 2] \cdot [\cos(2x) + \cos(2y)]. \quad (46)$$

The initial condition corresponds to an initial Mach number  $M \approx 0.08$ , with wavenumber  $k = 2\pi/\lambda = 1$ . Simulations were carried out on a hexahedral mesh of  $96^3$  and  $128^3$  resolution with 4th-order and 5th-order spatial discretisation schemes respectively. The WENO, CWENO, and CWENOZ variants were used with a CFL number of 1.3 for the explicit Runge-Kutta 4th-order scheme, up to  $t = 20$  for obtaining the dissipation statistics. The DNS

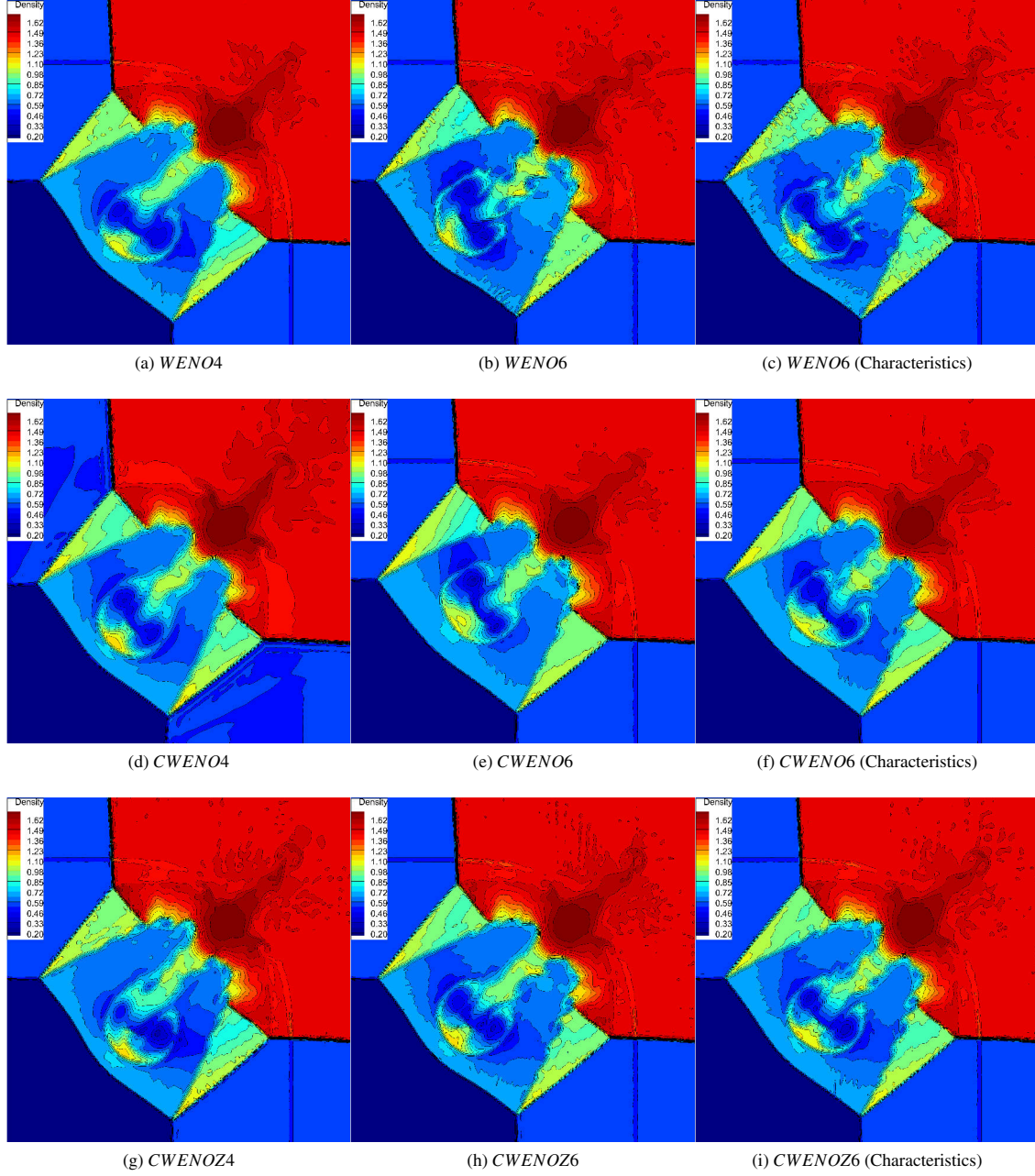


Figure 7: Density distribution with various 4th-order numerical schemes for the multidimensional Riemann problem at  $t_{final} = 1.0$ . Only the lower left portion of the computational domain is shown. It can be noticed that the all the schemes predict the correct solution pattern, but the WENO implementation provides significantly more oscillatory results and a more pronounced loss of symmetry compared to the other implementations.

results of Brachet et. al [111] are used for comparisons against the computed solutions. For numerical results for the 3D Taylor-Green vortex obtained with a new family of staggered semi-implicit discontinuous Galerkin finite element



schemes, see [112–115].

From the obtained results as shown in Fig. 9 it can be seen that the energy dissipation rate for all the 4th-order schemes is within the  $Re = 800 - 1600$  DNS results range, as the we move to the 5th-order schemes with a finer mesh the agreement with the  $Re = 1600$  DNS results improves as expected. The structure predicted near  $t \approx 5 - 6$  for the 5th-order scheme are due to the specific aspects of the ILES framework of the schemes, and are present in multidimensional numerical methods as previously reported in the previous studies by Drikakis et al. [106]. The key differences between the CWENO and CWENOZ schemes is a slightly larger secondary peak at  $t = 9.5$  for the  $96^3$  resolution, and an earlier peak for the CWENOZ scheme for the  $128^3$  resolution, whereas the results from the WENO schemes indicate a more dissipative behaviour for both settings.

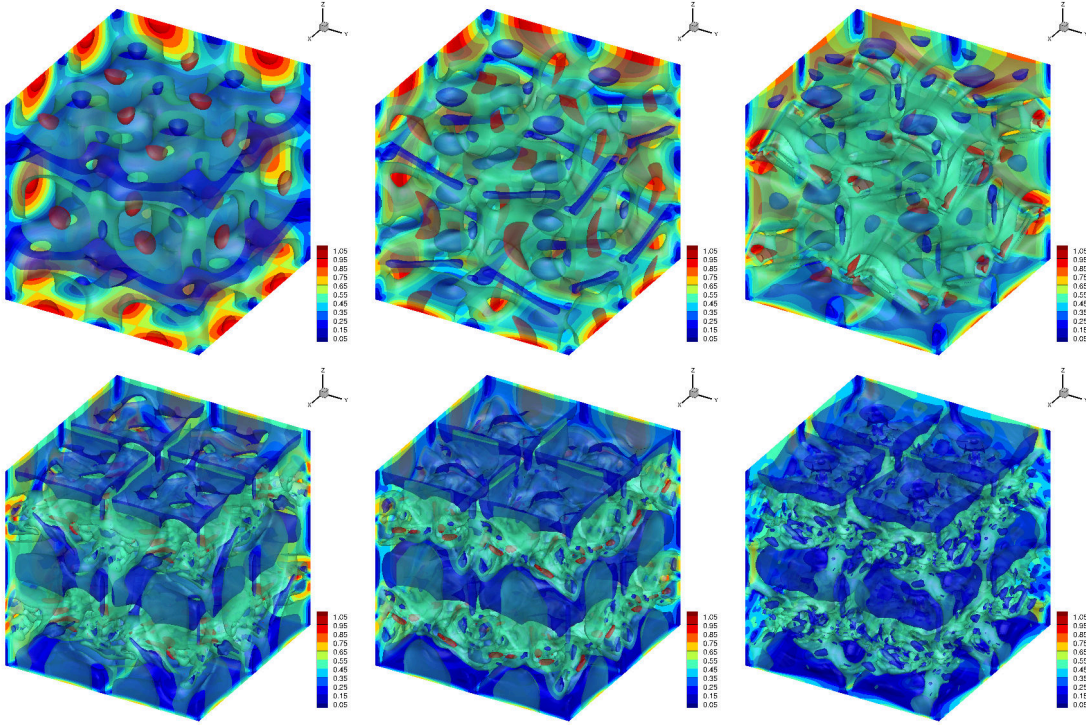


Figure 8: solution of the Taylor-Green vortex flow at  $Re=1600$  computed with the CWENO5 scheme on a hexahedral mesh of  $128^3$ . The isosurfaces of the velocity magnitude at three levels (0.1, 0.5, 1.0) are plotted at times  $t = 1.3, 2.7, 4.1, 6.4, 8.2$  and  $11$  from top left to bottom right respectively.

This test problem has been used to assess the computational costs of all the schemes. Ultimately we want to understand the computational savings that the CWENO and CWENOZ schemes can offer compared to the traditional WENO schemes. It needs to be stressed that the key memory requirement for each scheme is dependent of the storage of the Moore-Penrose pseudo-inverse of the matrix  $A_{mk}$  of Eq. (14), and this is where the key difference between a CWENO/CWENOZ and WENO schemes lie. For example for each hexahedral cell, for each conserved variable a WENO6 scheme would require approximately 338 kBytes of memory for storage of these matrices, while a CWENO/CWENOZ scheme require just 49kBytes of memory. Therefore significant savings in terms of storage are obtained with CWENO/CWENOZ reduced memory footprint. The communication of the values of for the halo cells arising from directional stencils become important (14% total time per iteration) at partitions of fewer than 160 elements per physical core for a 6th-order scheme as reported by Tsoutsanis et al. [86] for the present implementation of UCNS3D. Therefore the reduction of the size directional stencils for CWENO/CWENOZ schemes would prove beneficial for massively parallel simulations at partitions that include fewer than 160 elements per physical core.

The process involved running each of the schemes for 1000 iterations and averaging the total time taken per

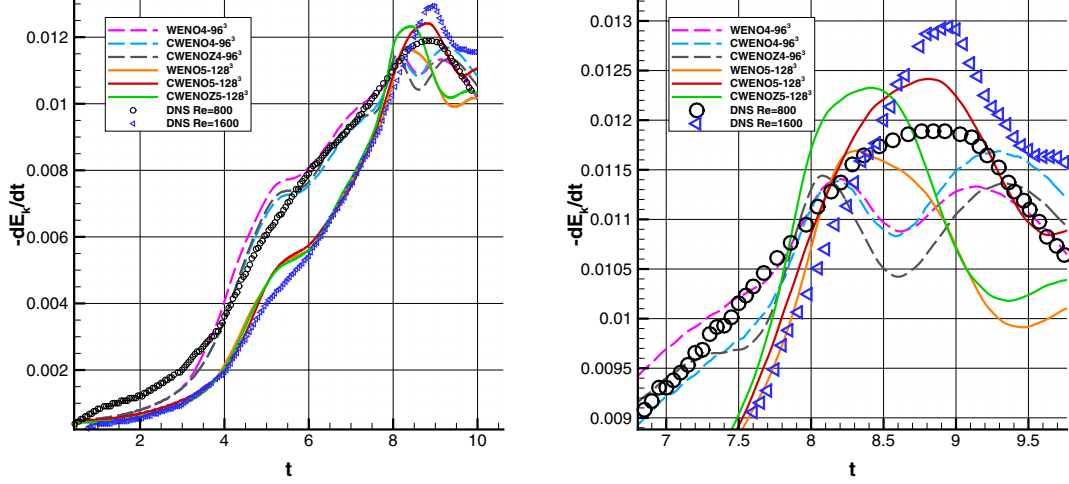


Figure 9: Kinetic energy dissipation rate for the viscous ILES of the Taylor-Green Vortex  $Re = 1600$  obtained with various schemes on a  $96^3$  and  $128^3$  hexahedral meshes, and comparison with the DNS results of Brachet et al. [111]. It can be noticed that all the schemes perform similarly, with CWENO and CWENOZ schemes being marginally less dissipative compared to WENO. The WENO4 scheme being 1.36 times more expensive computationally compared to CWENO and CWENOZ. Increasing the order of accuracy and grid resolution a better agreement with the DNS results is obtained.

iteration while excluding initialisation times, and I/O times. For this purpose, the University of Cambridge Tier-2 national high-performance computing service Peta4 was used, and in particular the Intel Xeon-Knights Landing cluster consisting of 342 nodes of Intel Xeon Phi 7210 with 96GB of RAM and Intel OPA at 100Gbs. All the simulations were executed with the utilisation of hybrid MPI+OPENMP implementation of the UCNS3D code, with 4 MPI processes per Knights Landing chip, and 64 threads spawned from each one of them, resulting a total of 256 threads per chip. This combination was found to be the ideal combination as thoroughly documented in [86]. All the simulations were executed in 20 Intel Xeon Phi 7210 processors, involving 80 MPI processes and 5,120 threads in total.

As expected the improvements in terms of computational times for the CWENO/CWENOZ schemes compared to WENO schemes become more prominent as the polynomial order is increased as seen in Table 6 at least for the reconstructions with respect to the conservative variables. On the other hand for the reconstruction in characteristic variables the computational cost is increased for all the schemes. But for the central schemes the computational cost increment is larger for WENO scheme when switching from conservative to characteristic variables. This attribute is linked to the additional computations required for the central schemes for the  $p_1$  polynomial according to Eq. (23), since the projection of the polynomial degrees of freedom to characteristic space for the calculation of the WENO weights is the most computational expensive process as reported in [86]. Therefore the characteristic variable reconstruction should be deployed only when really needed due to the increased computational cost.

Since we have not noticed significant differences in terms of accuracy between the current implementation of CWENO and CWENOZ schemes, but a slightly more robust non-oscillatory behaviour from the CWENO family, from this point onward the CWENO scheme will be utilised for the remaining test problems.

### 3.6. Schardin's Problem

The well established test problem of Schardin [116] is employed to assess the performance of the CWENO scheme, since this test problem involves multiple challenging flow structures interacting such as reflected shocks, Mach stems, contact waves, and vortices. A moving normal shock with Mach number  $M_s = 1.34$  is encountering a triangular wedge and the size of the computational domain is similar to the experimental one as seen in Fig. 10.

Table 6: Normalised computational times (with respect to CWENO3) for all the schemes for the viscous ILES of the Taylor-Green Vortex.

Polynomial Order	CWENO/Z (Conservative)	WENO (Conservative)	CWENO/Z (Characteristics)	WENO (Characteristics)
$\mathcal{P}2$	1.00	1.04	2.57	2.52
$\mathcal{P}3$	1.61	2.19	3.80	4.04
$\mathcal{P}4$	2.76	3.84	5.90	6.46
$\mathcal{P}5$	4.80	7.04	9.09	10.28

Table 7: Initial Conditions for Schardin's test problem

Parameter	Post-Shock	Pre-Shock
$P$ (Pa)	195557.25	101325
$T$ (K)	350.5	288.15
$\mathbf{u}$ (m/s)	(168.62,0,0)	(0,0,0)

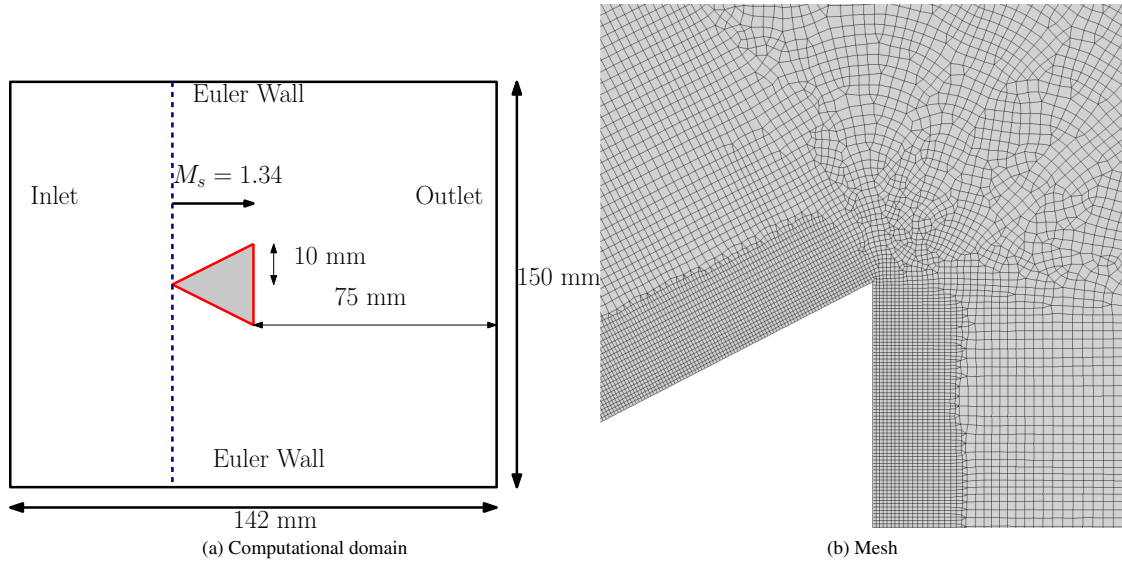


Figure 10: Computational domain with boundary conditions, and blown-up view of the mesh close to the triangular wedge for the Schardin's test problem.

The initial conditions are listed in Table 7, additionally the gas constant is  $R = 287 J/kg \cdot K$ , and the ratio of specific heats  $\gamma = 1.4$ . The hybrid computational mesh employed can be seen in Fig. 10 and consists of 5,264,594 cells with a resolution of  $h = 0.01 mm$  close to the wall, the prescribed boundary conditions correspond to a supersonic inflow and outflow, with inviscid wall boundary condition applied for the triangular wedge and the top and bottom sides of the domain as seen in Fig. 10. The inviscid Euler equations are solved with a CWENO5 scheme using characteristic variables reconstruction with a 4th-order (SSP) Runge-Kutta time stepping scheme while using  $CFL = 1.4$ . At the start time  $t = 0$  the incident shock wave contacts the triangular wedge, and the simulation is run for  $t = 200 ms$ . The obtained results are compared with the experimental imaging of Chang and Chang [117] at different instants as it can be seen in Fig. 11. From the results it can be noticed that all the flow features of the experiment are captured

at their correct location such as the bow shock, tip vortices, contact surfaces between the tip vortices and the shock triple point. The combination of the relatively fine mesh, and the CWENO5 is sufficient to capture the acoustic waves produced by the shock-vortexlet interactions as shown in Fig. 12, and agreeing with the experimental and computational results of Chang and Chang [117]. Finally the prediction of the Mach number along the symmetry line at the wake of the triangular wedge agrees well with the results of Chang and Chang [117] as seen in Fig. 12.

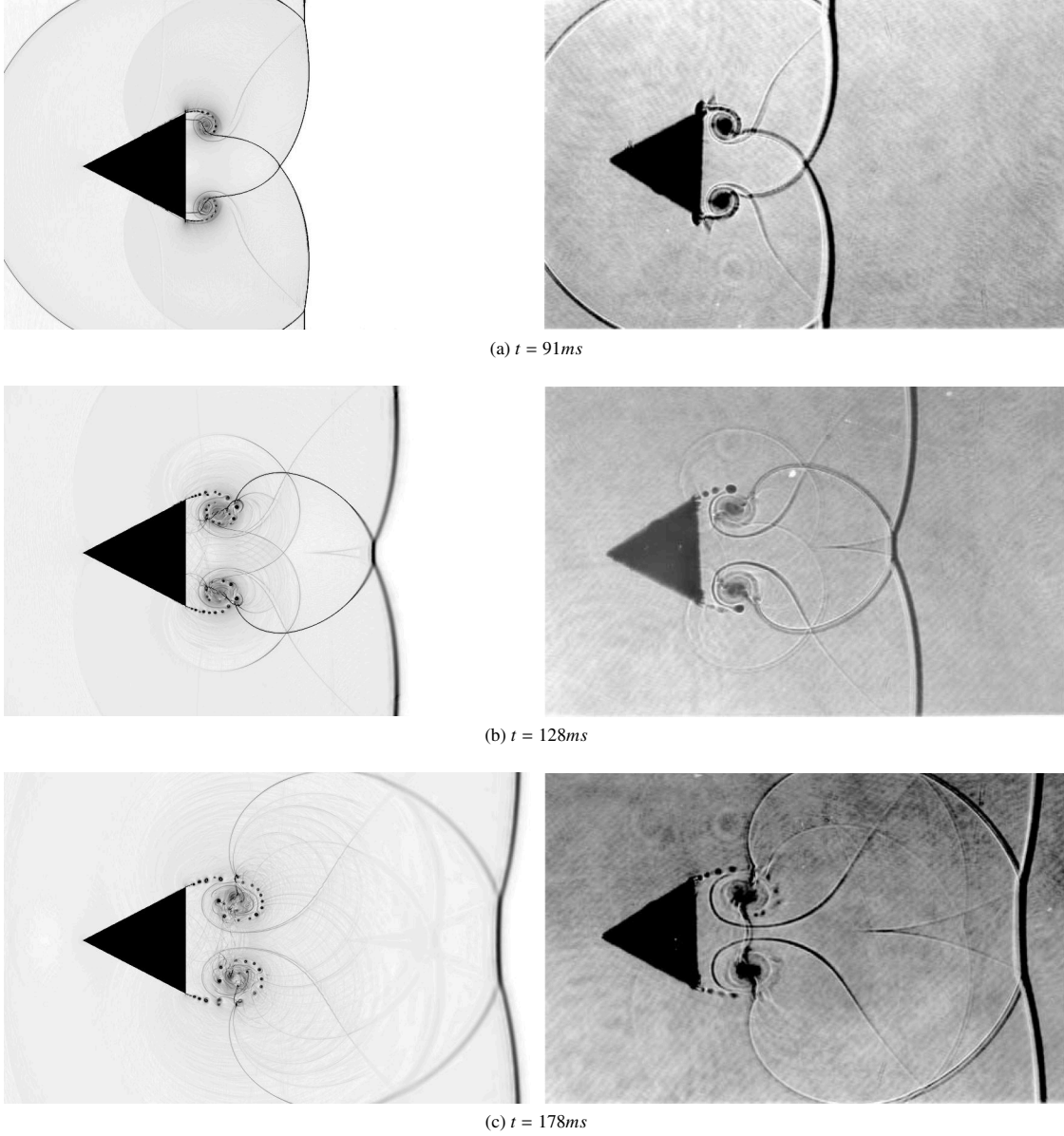


Figure 11: Contour plots of density gradient magnitude for the Schardin's test problem at various instants as computed with the present CWENO5 scheme (left), and comparison with the experimental results of Chang and Chang [117] (right)



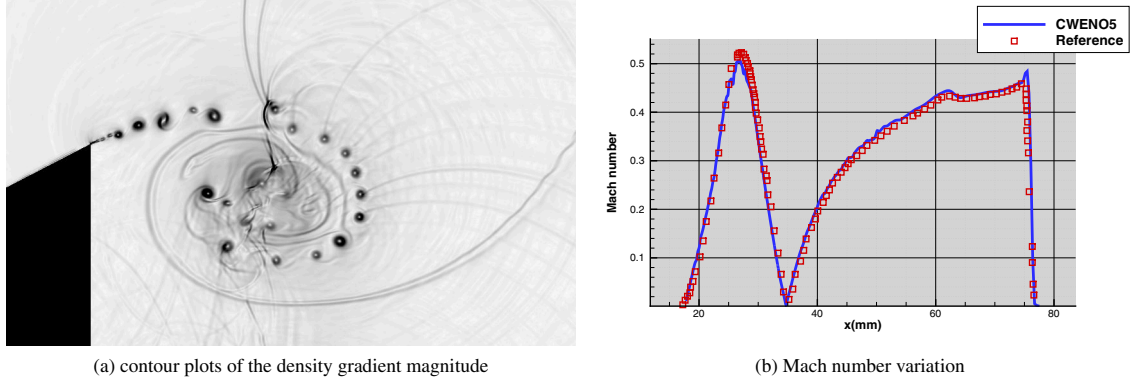


Figure 12: Contour plots of the density gradient magnitude for Schardin's test problem at the shock-vortexlet interaction zone at  $t = 172ms$  (left), and Mach number variation at the symmetry line wake of the triangular wedge and comparison with the results of Chang and Chang [117] (right).

### 3.7. ILES of Transonic cylinder $Re = 3900$

The last test case concerns the ILES of a transonic infinite cylinder of diameter  $D = 1$  at  $Re = 3900$  and  $M_\infty = 0.8$ . The majority of the available computational studies are focused on the subsonic ( $M_\infty = 0.2$ ) regime for this Reynolds number [29, 118–125] for subsonic Mach numbers. However, beyond  $M_\infty = 0.75$  a multitude of phenomena occur including shock wave boundary layer, shock wave shear layer and shocklet/turbulence interaction, forming an ideal environment for assessing the robustness and fidelity of the developed schemes.

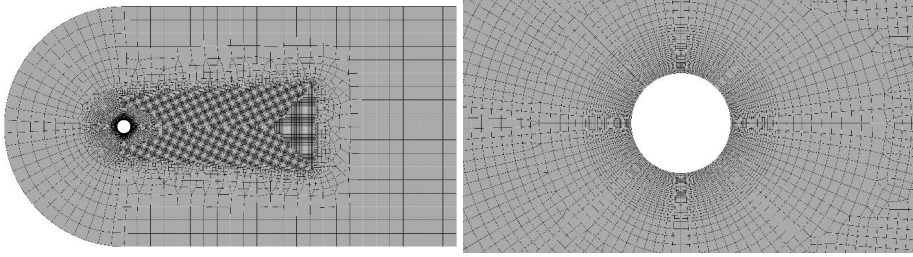


Figure 13: Overview of the hybrid unstructured mesh used for the transonic flow past an infinite cylinder at  $Re = 3900$  and  $M_\infty = 0.8$ .

For this case a very coarse hybrid unstructured mesh of approximately 0.8 million cells is used, consisting of hexahedral and prismatic cells as illustrated in Fig. 13. The domain extends  $10D$  upstream and  $25D$  downstream, and  $\pi \cdot D$  in the  $z$ -direction. The grid resolution at the boundary layer region gives a  $y^+ \approx 1$  at the first cell-centre off the surface, 64 cells are used in the  $z$ -direction, and a refinement region at the wake of the cylinder is used for  $10D$  downstream with an average cell length of 0.09. Periodic boundary conditions are used in the  $z$ -direction, no-slip boundary conditions at the surface of the cylinder and free-stream conditions at the farfield. The CWENO 4th-order scheme is used, with a CFL number of 1.45. The simulation was run for  $t = 50t_c$  where  $t_c = D/U_\infty$  for the initial transients to settle, and for an additional  $240t_c$  for obtaining sufficiently converged statistical data. The variation of lift and drag coefficient can be seen in Fig. 14, with the averaged drag coefficient from this study being within the range of results provided by previous computational studies, experiments and flight test as listed in Table 8 at the same Mach number, although the Reynolds number is not the same.

From the time-averaged and spanwise-averaged pressure coefficient obtained as seen in Fig. 15 the results obtained are in reasonable agreement with the computational results of Xia et al. [126]  $Re = 4 \times 10^4$  (CLES) in terms of pressure coefficient up until  $70^\circ$ . Beyond that the separation and reattachment occur at different locations as seen from the skin friction coefficient, which is mainly attributed to the different Reynolds number. From the instantaneous flow-field

Table 8: Comparison of drag coefficient with experimental and computational results at Mach=0.8

Source	$C_D$
Present CWENO4 (ILES)	1.408
Xia et al.[126] $Re = 4 \times 10^4$ (CLES)	1.353
Xia et al.[126] $Re = 1 \times 10^6$ (CLES)	1.312
Welsh [127] (Flight Test)	1.425
Macha [128] (Experiment)	1.529

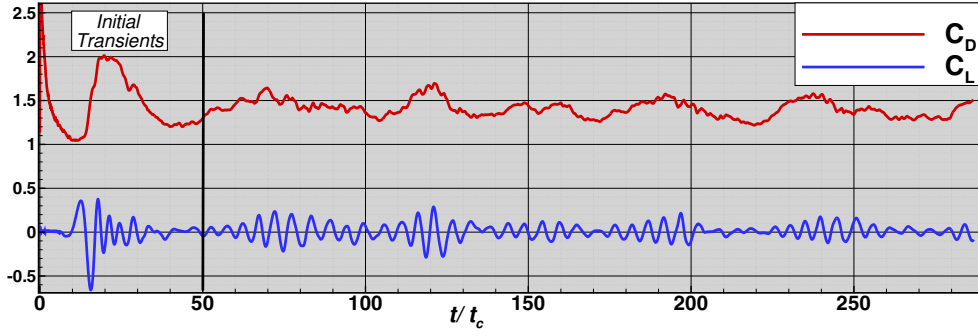


Figure 14: Lift and drag coefficients variations in time for the flow past a cylinder at  $Re = 3900$  and  $M_\infty = 0.8$ .

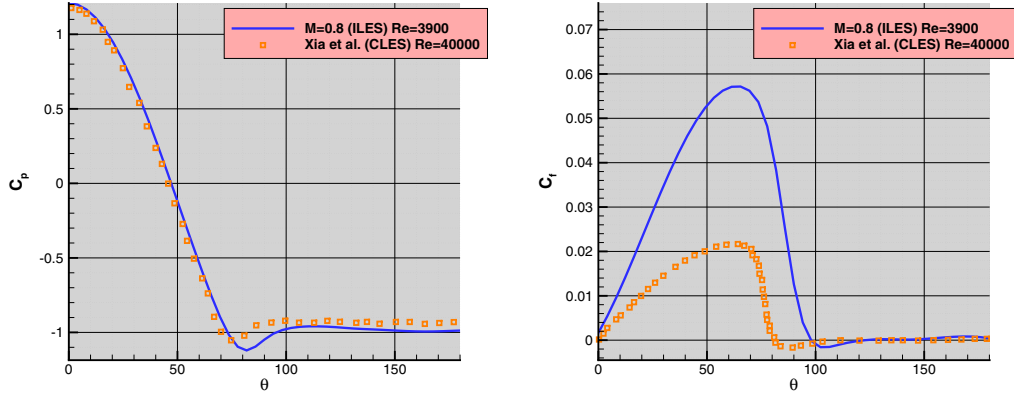


Figure 15: Time-averaged and spanwise-averaged pressure coefficient and skin friction coefficient for the flow past a cylinder at  $Re = 3900$  and  $M_\infty = 0.8$  and comparison with the results of Xia et al.[126]  $Re = 4 \times 10^4$  and  $M_\infty = 0.8$  (CLES).

solution as seen in Fig. 16 it can be noticed that all the flow features typical of this regime such as the shocks at the shoulders of the cylinder, and the long shear layers interaction with the shocklets at the near wake agree with previous studies of Xia et al. [126] .

Finally, by placing probes at different locations in the wake of the cylinder the dominant vortex shedding frequency was found to be  $St = 0.185$  as seen in Fig. 17, which is lower compared to the subsonic case of  $M_\infty = 0.2$  for the same Reynolds number as reported in [29, 118–125], but agrees well with the trends noticed in previous experimental

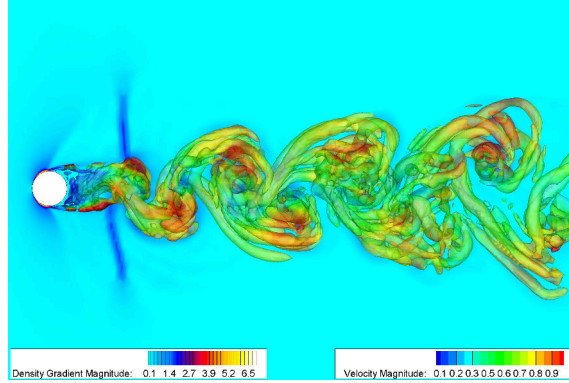


Figure 16: Isosurfaces of Q-criterion coloured by velocity magnitude, and contour plots of the density gradient magnitude on a 2D slice at  $z=0$ , for the flow past a cylinder at  $Re = 3900$  and  $M_\infty = 0.8$ , where the interaction of the vortical structures with shock waves typical of this flow regime can be noticed.

and computational studies as reported by Canuto and Taira [129] where the Strouhal number is reduced when Mach number increases.

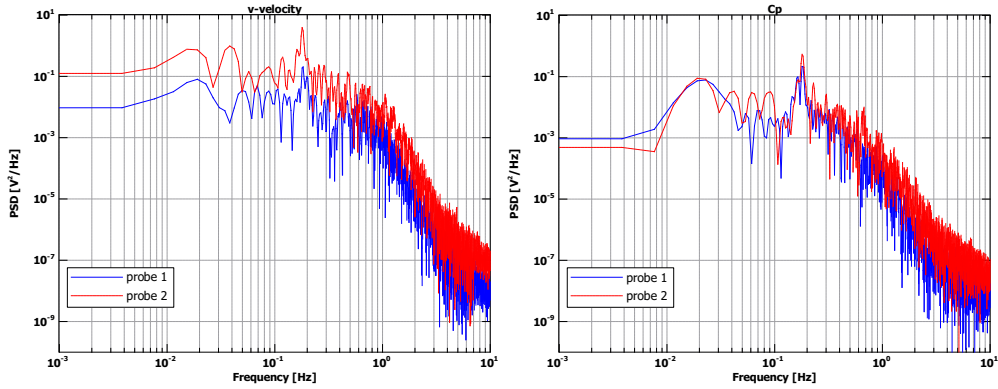


Figure 17: 1D frequency spectra for two probe positions located at  $(1.30, 0.69, \pi/2)$  and  $(2.0, 0.59, \pi/2)$  the transonic flow past an infinite cylinder at  $Re = 3900$  and  $M_\infty = 0.8$ .

#### 4. Conclusions

This paper extends the central high-order non-oscillatory family of schemes to mixed-element unstructured meshes up to 6th-order of spatial accuracy. The performance benefits obtained from the CWENO and CWENOZ class of schemes are not only related to a decreased computational time compared to the WENO schemes but significantly improved robustness. The latter being linked with the compactness of the directional stencils that increase the probability of at least one of them lying in smooth-data regions. The CWENO scheme exhibits an increased sensitivity compared to CWENOZ and WENO with respect to the value of the linear weight assigned to the central stencil, since a large value is required to achieve the designed order of accuracy for mixed-element unstructured meshes. On the other hand the CWENO demonstrated the most robust non-oscillatory behaviour in all the test problems, and was therefore the method of choice for the last two test problems which included a plethora of smooth and discontinuous flow features, where a good agreement with experimental results was obtained.

The computational cost of the new class of schemes can be approximately up to 30% less than the equivalent WENO scheme in 3D when conservative variables reconstruction are used. However, when switching to characteristic variables the cost of the schemes is approximately up to 11% less compared to WENO schemes. For characteristic variables reconstruction the most expensive procedure is the projection of the polynomial degrees of freedom to characteristic space. For the CWENO/CWENOZ schemes the additional operations required by the  $p_1$  polynomial are significantly greater compared to the central stencil polynomial of the WENO schemes. The latter is not considered an important drawback due to the reduced complexity of the algorithm, and the reduced size of the directional stencils that will translate to reduced communications across processes and reduced memory footprint for massively parallel computations.

Expanding the current family of schemes to the targeted ENO schemes [130–132] for further reducing their computational footprint is an interesting line of future research. Future developments will concern the extension of the subject schemes to overset meshes for industrial-scale applications since the compact nature of the present non-oscillatory schemes can be useful for interpolating the solutions across different meshes. Future applications of the method proposed in this paper will concern the unified first order hyperbolic model of continuum mechanics proposed and discussed in [59, 133, 134], as well as the novel first order hyperbolic reformulations of nonlinear dispersive systems forwarded in [135, 136]. With the high order CWENOZ schemes developed in this paper we also plan to study natural convection problems in the low Mach number limit, similar to the problems investigated in [137, 138].

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## **Arbitrary High Order Central Non-Oscillatory Schemes on Mixed-Element Unstructured Meshes**

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# Arbitrary high order central non-oscillatory schemes on mixed-element unstructured meshes

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